

A New Era in High Speed Analyses for Food and Environmental Applications

(Agilent 1290 Infinity uHPLC,
Agilent 6230 TOF, 6460 QQQ and 6540 Q-TOF.)

Peter Stone
Agilent Technologies Inc,
Santa Clara, CA.



Agilent Technologies

Agilent 1290 Infinity LC Attributes for MS

- *Infinitely Better for LC/MS*



Lowest
Delay
Volume

- Pump (w/o) mixer: 10 μ L
- Pump, Fixed Loop 20 μ L
- Pump, Fixed Loop, JetWeaver 55 μ L

Highest
Precision

- ALS precision for small volumes:
 <1.5% from 0.5-1 μ Lm, <0.7% from 1-2 μ L,
 <0.25% @ 2-20 μ l (40 μ l) **
- Pump Active Damping:
 RT stability < 0.2 % (1.5 min runs)**

Best
Autosampler
Performance

- <0.002% carry-over with Chlorhexidine
- Optional needle seat backflushing with FlexCube
- Fixed Loop or Variable Loop Injections

Greatest
Productivity

- 1200bar @ 2mL/min for highest resolution per time
 Reduced Ion & Matrix Suppression
- HT-Solution for up to 2000 samples/day (ACR)
- Complete Integration and control from MassHunter
- Enables method conversion from/to any (U)HPLC



Agilent Technologies

Fast Target Screening of PPCPs Using high sensitivity 6460 Triple Quadrupole with 1290 uHPLC

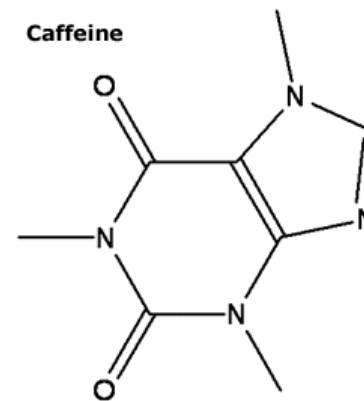
- ❑ Subfemtogram, on-column sensitivity
- ❑ Five Orders of Linearity
- ❑ 3,000 m/z mass range
- ❑ Fast Pos/Neg switching, 30 millisec
- ❑ New Method Development Tool – MH Optimizer
- ❑ “**Dynamic Multiple Reaction Monitoring**” for MRMs scheduled by peak retention times instead of time segments



Target Screening of PPCPs in Water Matrices LC/MS/MS QQQ

Pharmaceuticals and Personal Care Products (PPCPs) refer, in general, to any product used by individuals for personal health or cosmetic reasons or used by agribusiness to enhance growth or health of livestock.

PPCPs comprise a diverse collection of thousands of chemical substances, including prescription and over-the-counter therapeutic drugs, veterinary drugs, fragrances, and cosmetics.



EPA 1694 - Identification of PPCPs in Water

Antibiotics

Acetaminophen
Albuterol
Ampicillin
Anhydrochlortetracycline (ACTC)
Anhydrotetracycline (ATC)
Azithromycin
Caffeine
Carbadox
Carbamazepine
Cefotaxime
Chlortetracycline (CTC)
Cimetidine
Ciprofloxacin
Clarithromycin
Clinafloxacin
Cloxacillin
Codeine
Cotinine
Dehydronifedipine
Demeclocycline
Digoxigenin
Digoxin
Diltiazem
1,7-Dimethylxanthine
Diphenhydramine
Doxycycline
Enrofloxacin
4-Epianhydrochlortetracycline
(EACTC)
4-Epianhydrotetracycline (EATC)
4-Epichlortetracycline (ECTC)
4-Epioxytetracycline (EOTC)
4-Epitetracycline (ETC)
Erythromycin
Erythromycin anhydride
Flumequine
Fluoxetine
Gemfibrozil

Anti-Depressants

Pain Killers

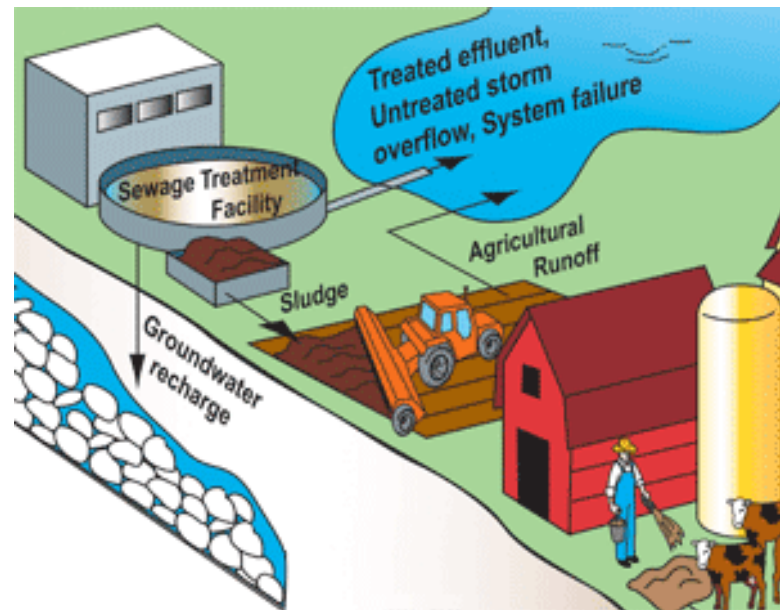
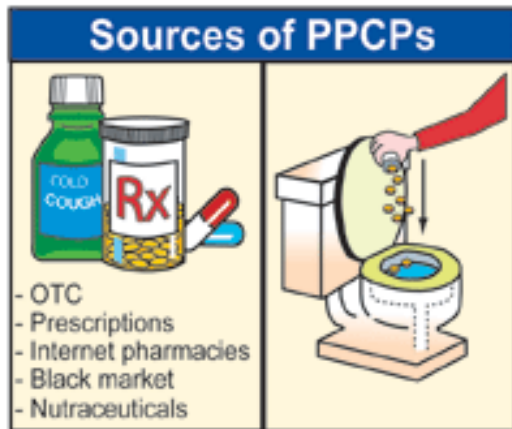
Ibuprofen
Isochlortetracycline (ICTC)
Lincomycin
Lofexidine
Metformin
Miconazole
Minocycline
Naproxen
Norfloxacin
Norgestimate
Ofloxacin
Ormetoprim
Oxacillin
Oxolinic acid
Oxytetracycline (OTC)
Penicillin V
Penicillin G
Ranitidine
Roxithromycin
Sarafloxacin
Sulfachloropyridazine
Sulfadiazine
Sulfadimethoxine
Sulfamerazine
Sulfamethazine
Sulfamethizole
Sulfamethoxazole
Sulfanilamide
Sulfathiazole
Tetracycline (TC)
Thiabendazole
Triclocarban
Triclosan
Trimethoprim
Tylosin
Virginiamycin
Warfarin
Other standards

Birth Control Steroids



Where are PPCPs found in the environment?

PPCPs in the environment are frequently found in aquatic environments because PPCPs dissolve easily and don't evaporate at normal temperature and pressure. Practices such as the use of sewage sludge ("biosolids") and reclaimed water for irrigation bring PPCPs into contact with the soil (landfill also).



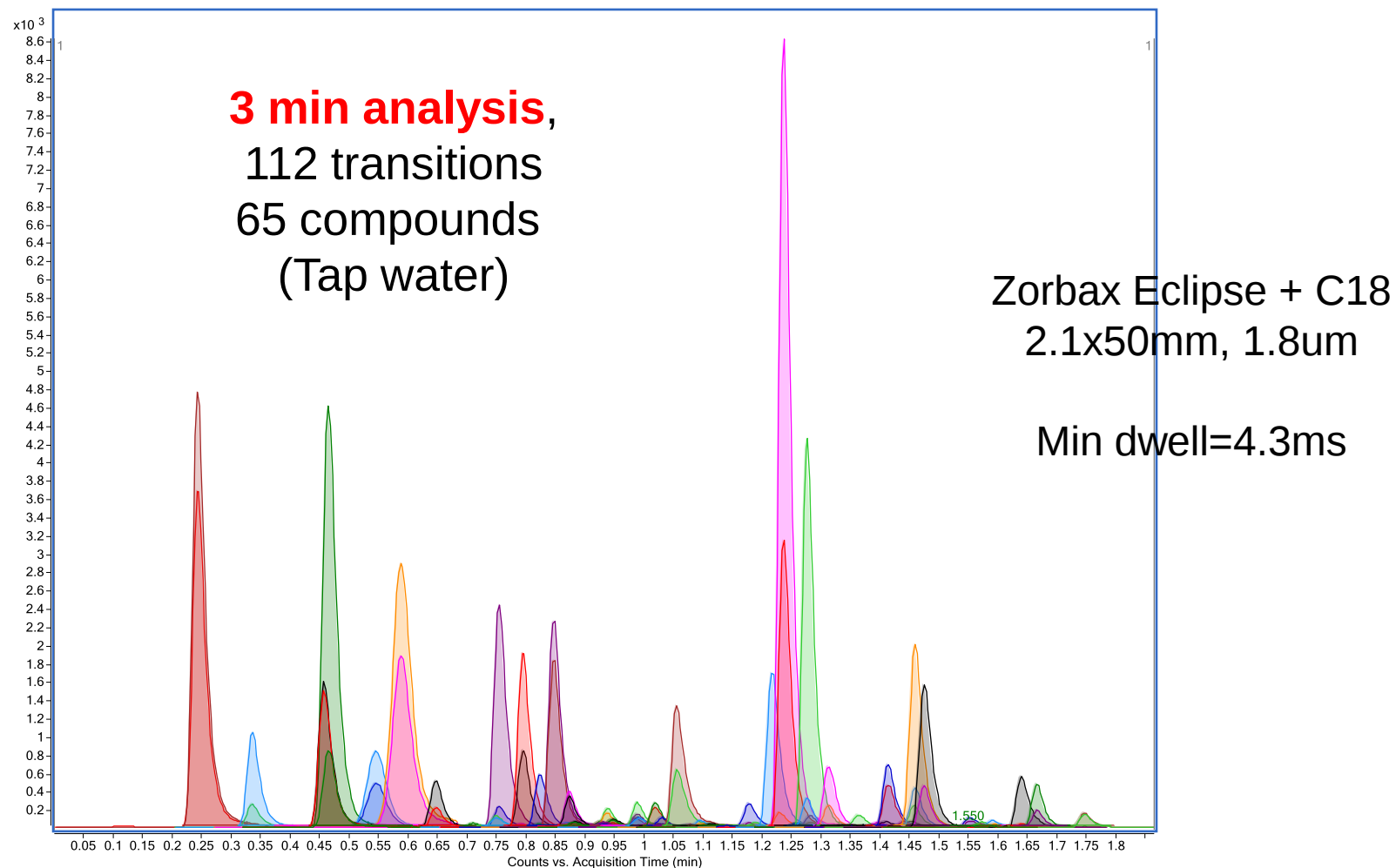
List of PPCP Target Analytes (EPA 1694, Positive Analysis)

Compound Name	Precursor Ion	Product Ion	Fragmentor (V)	Collision Energy (V)	RT	RT Window	Ion Polarity
1,7-Dimethylxanthine	181.0	124.0	90	15	0.658	0.50	Positive
4-Epi-anhydrotetracycline	427.0	410.0	90	15	3.473	0.50	Positive
4-Epichlortetracycline	479.0	462.0	134	17	2.416	0.50	Positive
4-Epitetracycline	445.0	410.0	110	15	1.777	0.50	Positive
Acetaminophen	152.0	110.0	90	15	0.656	0.50	Positive
Albuterol	240.0	148.0	90	15	0.489	0.50	Positive
Amphetamine	136.1	91.0	70	13	0.880	0.50	Positive
Anhydrotetracycline	427.0	410.0	90	15	3.473	0.50	Positive
Atenolol	267.2	145.0	134	21	0.502	0.50	Positive
Azithromycin	749.5	591.4	130	30	3.254	0.50	Positive
Caffeine	195.0	138.0	110	15	1.072	0.50	Positive
Carbadox	263.0	130.0	80	35	3.097	0.50	Positive
Chlorotetracycline	479.0	462.0	110	15	2.416	0.50	Positive
Cimetidine	253.0	159.0	100	10	0.489	0.50	Positive
Clarithromycin	748.5	158.0	110	25	4.384	0.50	Positive
Clonidine	230.0	44.0	150	25	0.764	0.50	Positive
Cloxacillin degradate	469.0	160.0	70	15	4.625	0.50	Positive
Cloxacillin	436.0	160.0	90	15	4.557	0.50	Positive
Codeine	300.0	215.0	130	25	0.709	0.50	Positive
Cotinine	177.0	80.0	90	25	0.420	0.50	Positive
Dehydronifedipine	345.0	284.0	130	25	4.662	0.50	Positive
Demeclocycline	465.0	448.0	130	15	2.394	0.50	Positive
Dextromethorphan	272.2	170.9	152	41	3.365	0.50	Positive
Diazepam	285.1	193.0	162	33	4.818	0.50	Positive
Digoxigenin	391.0	355.0	90	15	3.036	0.50	Positive
Diltiazem	415.0	178.0	130	25	3.702	0.50	Positive
Diphenhydramine	256.0	167.0	70	15	3.463	0.50	Positive
Doxycycline	445.0	428.0	110	15	3.008	0.50	Positive
Enrofloxacin	360.0	316.0	130	15	2.237	0.50	Positive
Erythromycin Anhydrate	716.5	158.0	90	25	4.201	0.50	Positive
Erythromycin	734.5	158.0	90	35	3.854	0.50	Positive
Erythromycin-Labeled	736.5	578.0	90	15	3.851	0.50	Positive

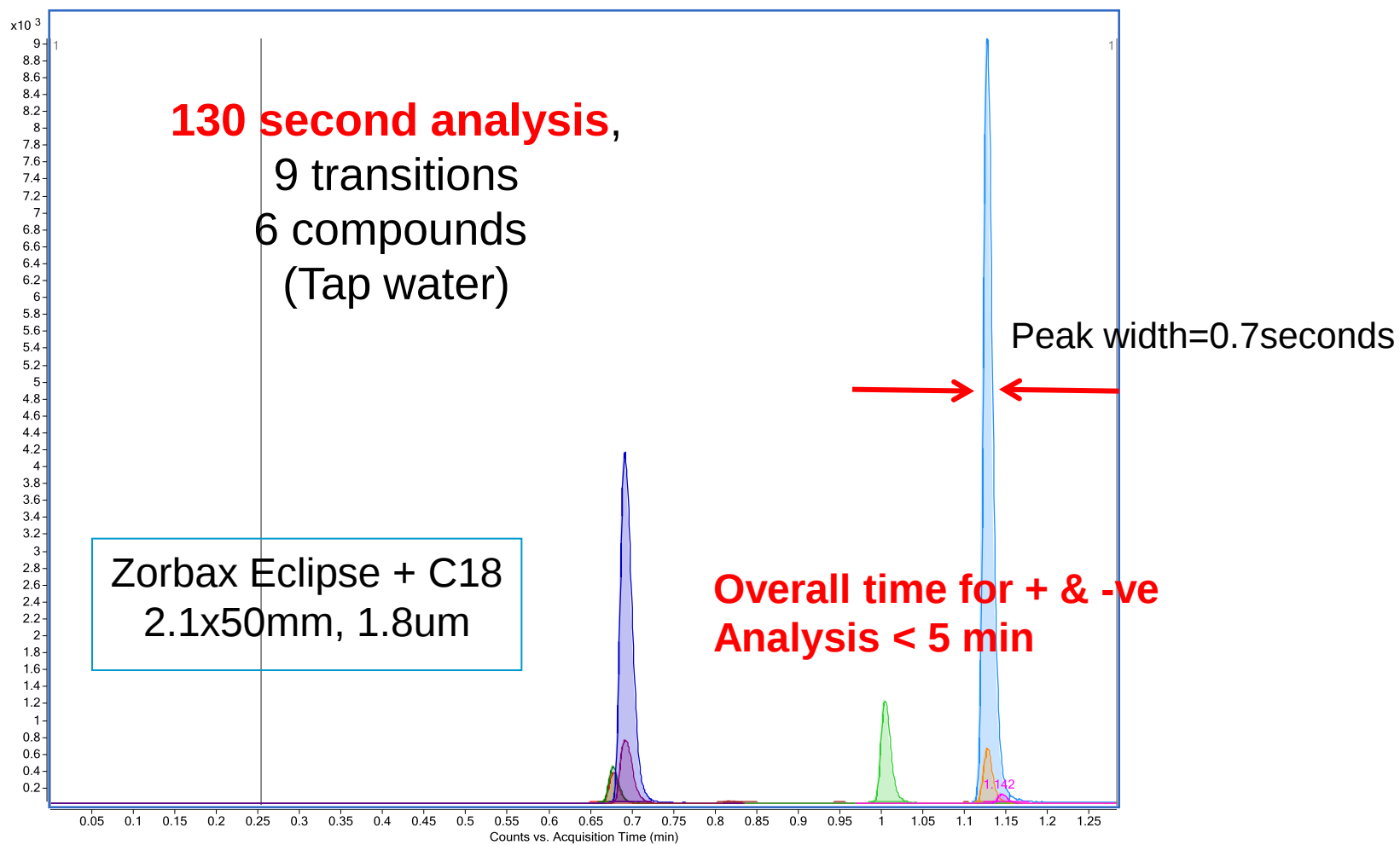
33/65 PPCP Analytes (Only quantifier ion transitions shown)



uHPLC - PPCPs from EPA 1694 (**Positive Polarity**)

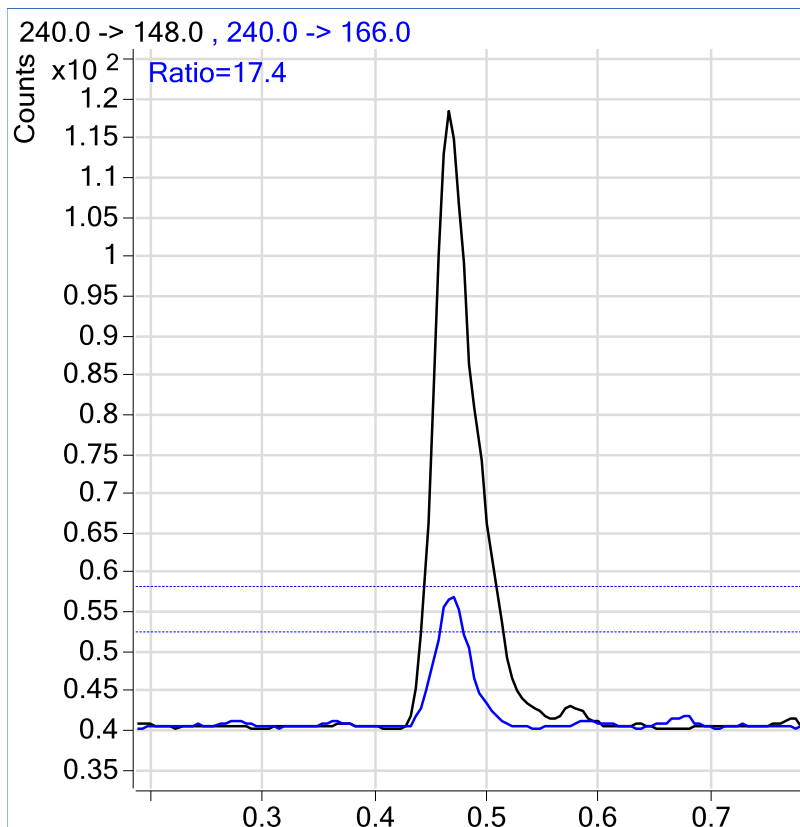


uHPLC - PPCPs from EPA 1694 (Negative Polarity)



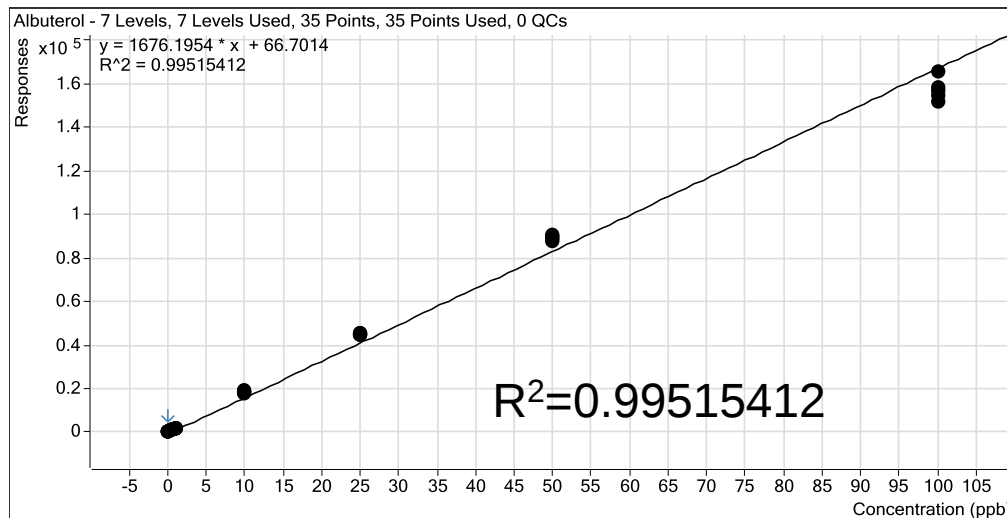
Chromatographic Precision (1290/6460, dynamic MRM)

Albuterol, 100fg on-column Tap Water Spike



Quant & Qual transitions
20% outlier ratio boundary

N= 5 (@100fg), %RSD (area) = 5.436907
N= 35 (all Cal levels), %RSD (RT) = 0.5182964

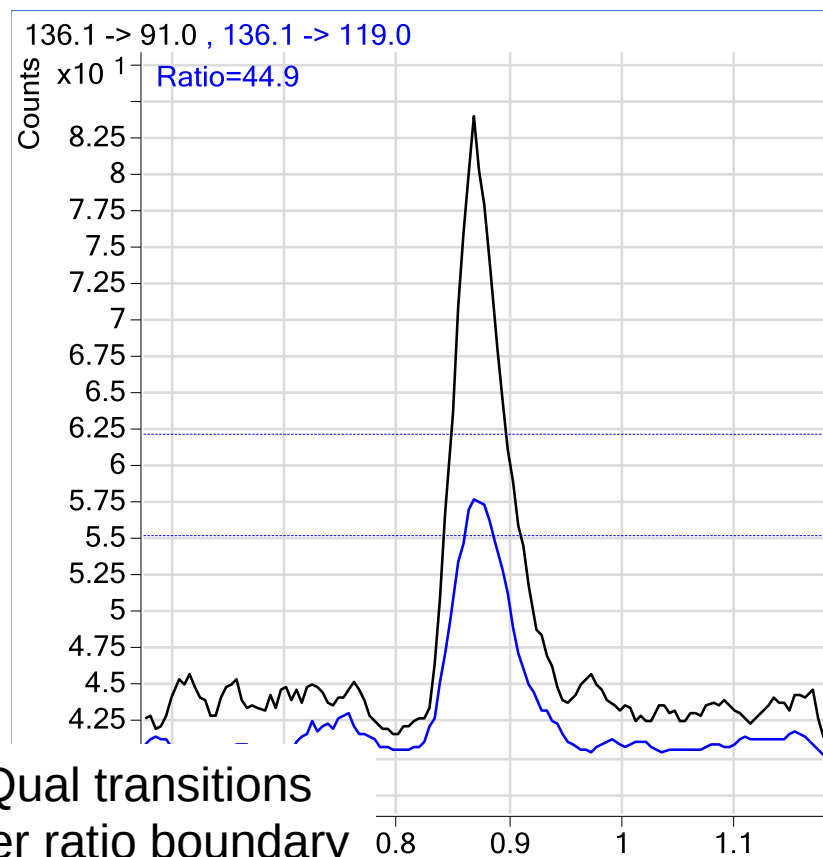
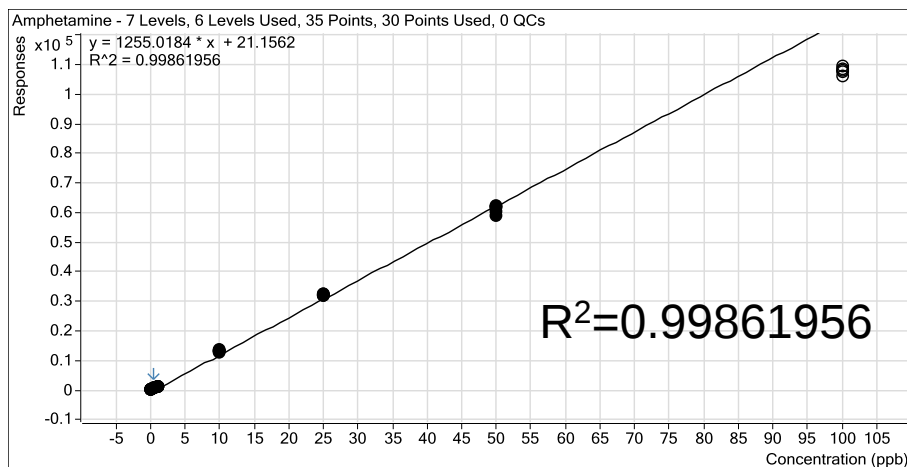


Chromatographic Precision (1290/6460, dynamic MRM)

Amphetamine, 100fg on-column Tap water spike

N= 5 (@100fg), %RSD (area) = 6.902997

N= 35 (all Cal levels), %RSD (RT) = 0.50386361

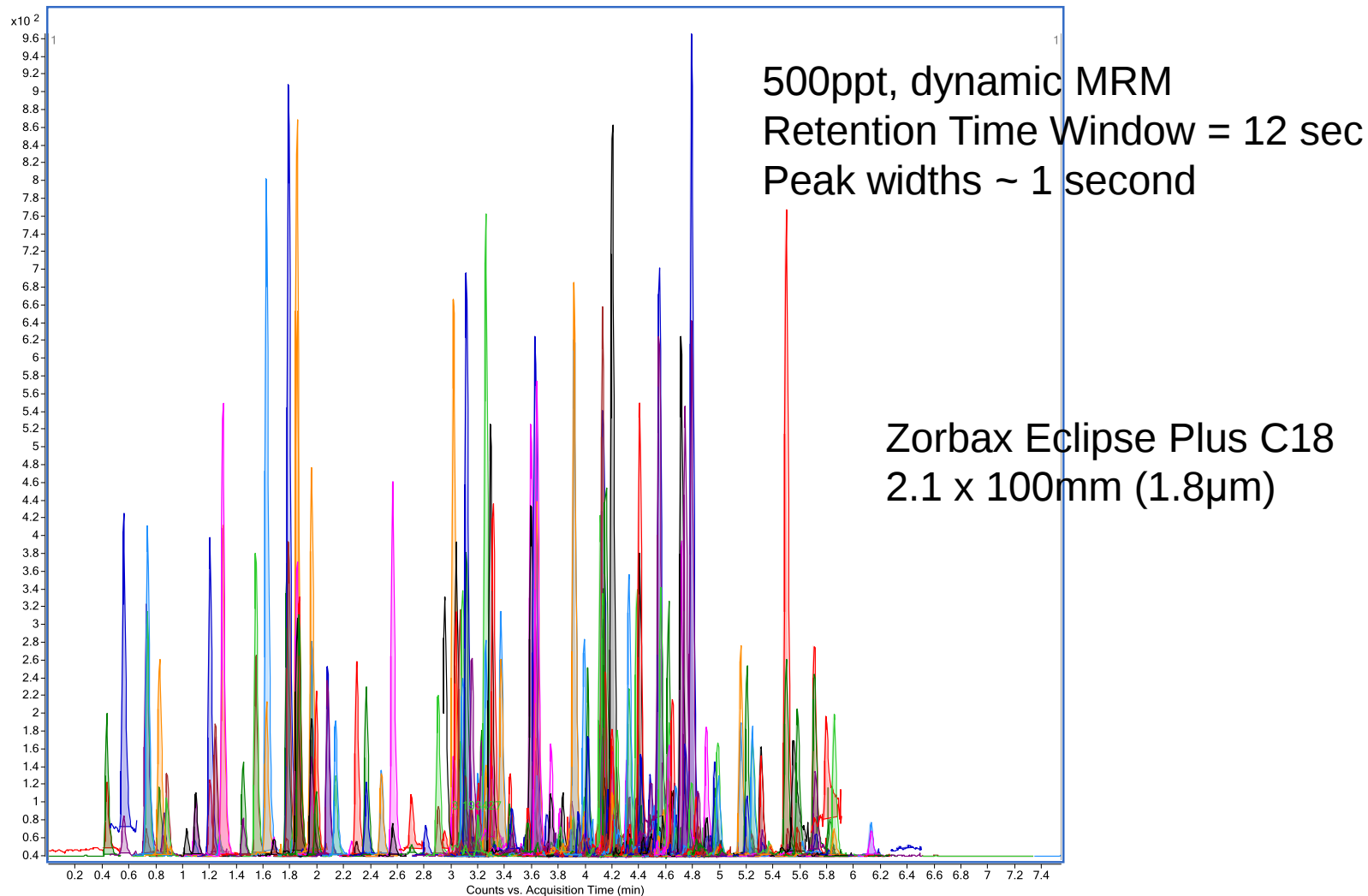


Quant & Qual transitions
20% outlier ratio boundary



8 Minute Dynamic MRM Analysis 6460 QqQ

- 250 Pesticide Screen.



6460 QqQ Method Conditions

MS QQQ

Acquisition **Source** Chromatogram Instrument Diagnostics

Source parameters

Gas Temp: 225 °C 287 °C

Gas Flow: 8 l/min 3.0 l/min

Nebulizer: 45 psi 15.0 psi

Sheath Gas Temp: 370 °C 178 °C

Sheath Gas Flow: 11 l/min 11.0 l/min

Positive Negative

Capillary: 4500 V 3500 V 7 nA

Nozzle Voltage: 500 V 1500 V

Chamber Current 0.23 µA

Copy

Paste

Paste to All Segments

Acquisition Source Chromatogram Instrument Diagnostics

Scan segments

Compound Name	ISTD?	Precursor Ion	MS1 Res	Product Ion	MS2 Res	Fragmentor	Collision Energy	Ret Time (min)	Delta Ret Time	Polarity
▶ Methamidophos	☐	142	Unit	125	Unit	80	10	0.435	0.2	Positive
Methamidophos	☐	142	Unit	94	Unit	80	10	0.435	0.2	Positive
Acephate	☐	184	Unit	143	Unit	90	5	0.557	0.2	Positive
Acephate	☐	184	Unit	95	Unit	90	20	0.557	0.2	Positive
Formetanate	☐	222.1	Unit	165.1	Unit	100	10	0.724	0.2	Positive
Formetanate	☐	222.1	Unit	120.1	Unit	100	25	0.724	0.2	Positive
Omethoate	☐	214	Unit	183	Unit	90	5	0.735	0.2	Positive
Omethoate	☐	214	Unit	125	Unit	90	20	0.735	0.2	Positive
Propamocarb	☐	189	Unit	101.9	Unit	100	15	0.816	0.2	Positive

Dynamic MRM Parameters

Cycle Time 300 ms

Total MRMs = 438 Max Concurrent MRMs = 52 Min/Max Dwell = 2.27 ms/296.50 ms



MassHunter “Dynamic MRM”

Many applications require screening of 100 – 1000 compounds in one MRM method !

- Food, environmental and Forensic analysis (e.g. Doping Control)
- Subsequent quantitation of positives in dedicated methods

WITHOUT Dynamic MRM

- Need to manually set up multiple time segments to maximize dwell time
- Tedious and fragile for chromatographic time shifts

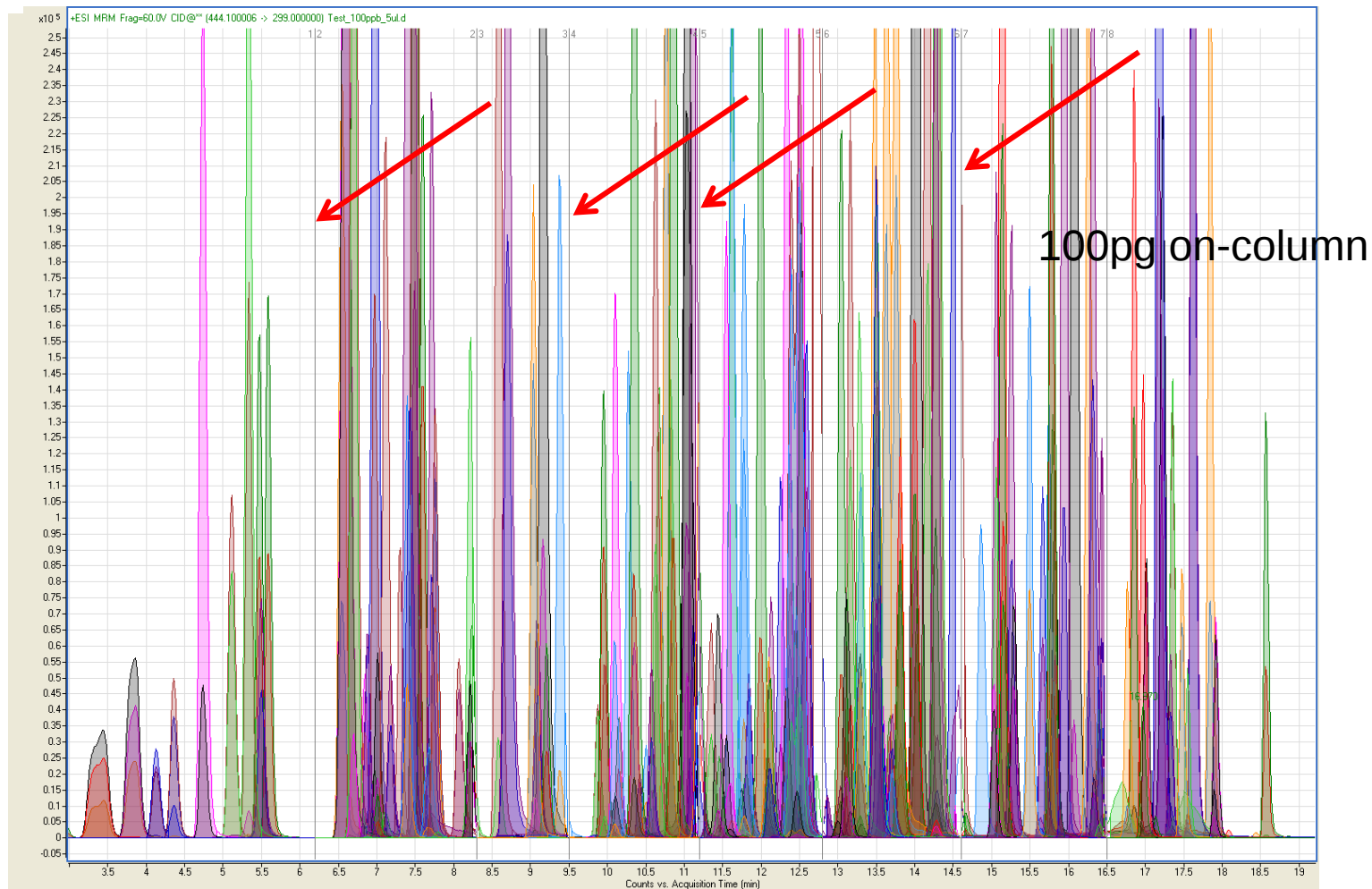
WITH Dynamic MRM

- Automatic setup of overlapping time segments without user intervention
- Even less MRMs per time results in even longer dwell time / sensitivity
- Unaffected by chromatographic time shifts

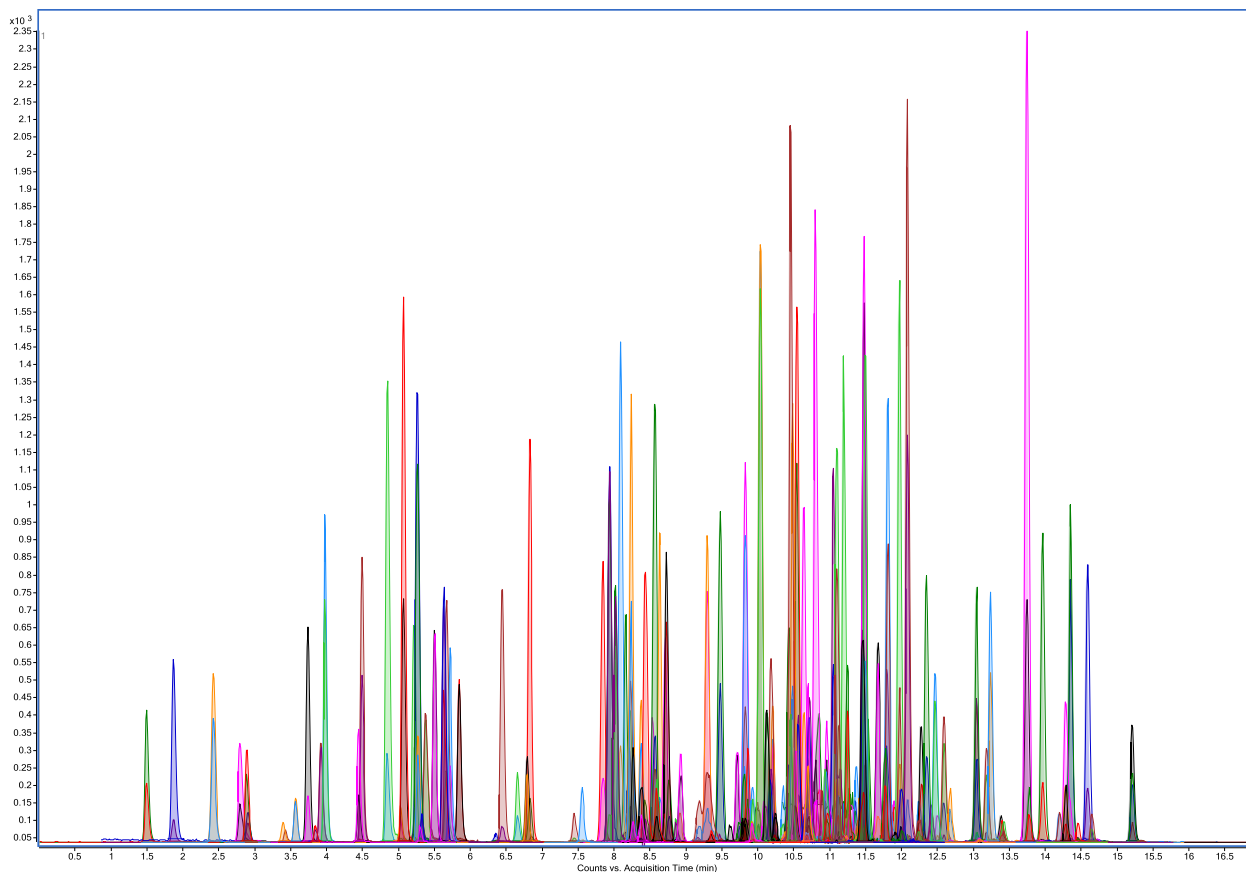


Segmentation with NO Dynamic MRM functionality

300 Component Standard



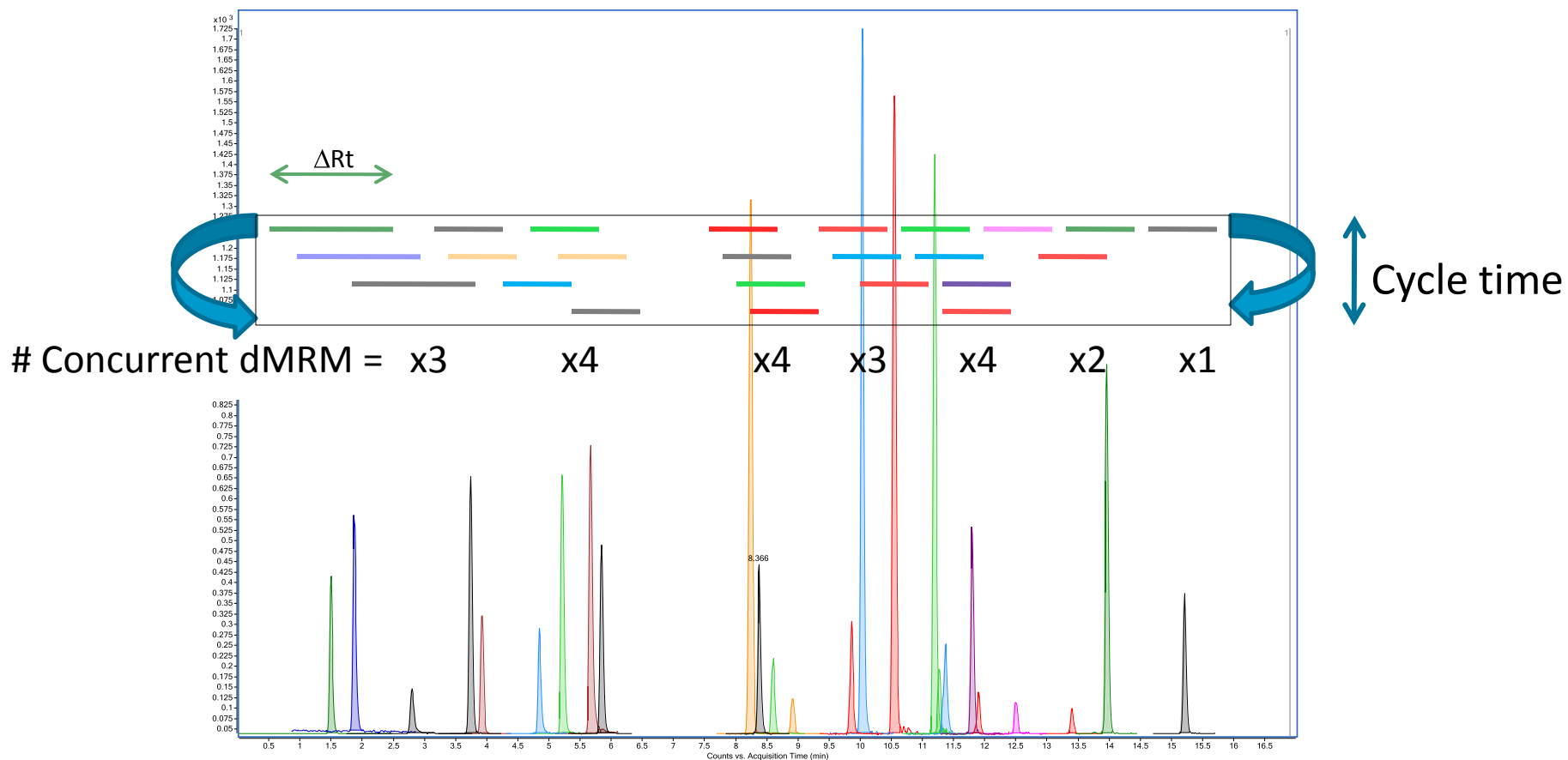
NO Segmentation with Dynamic MRMs



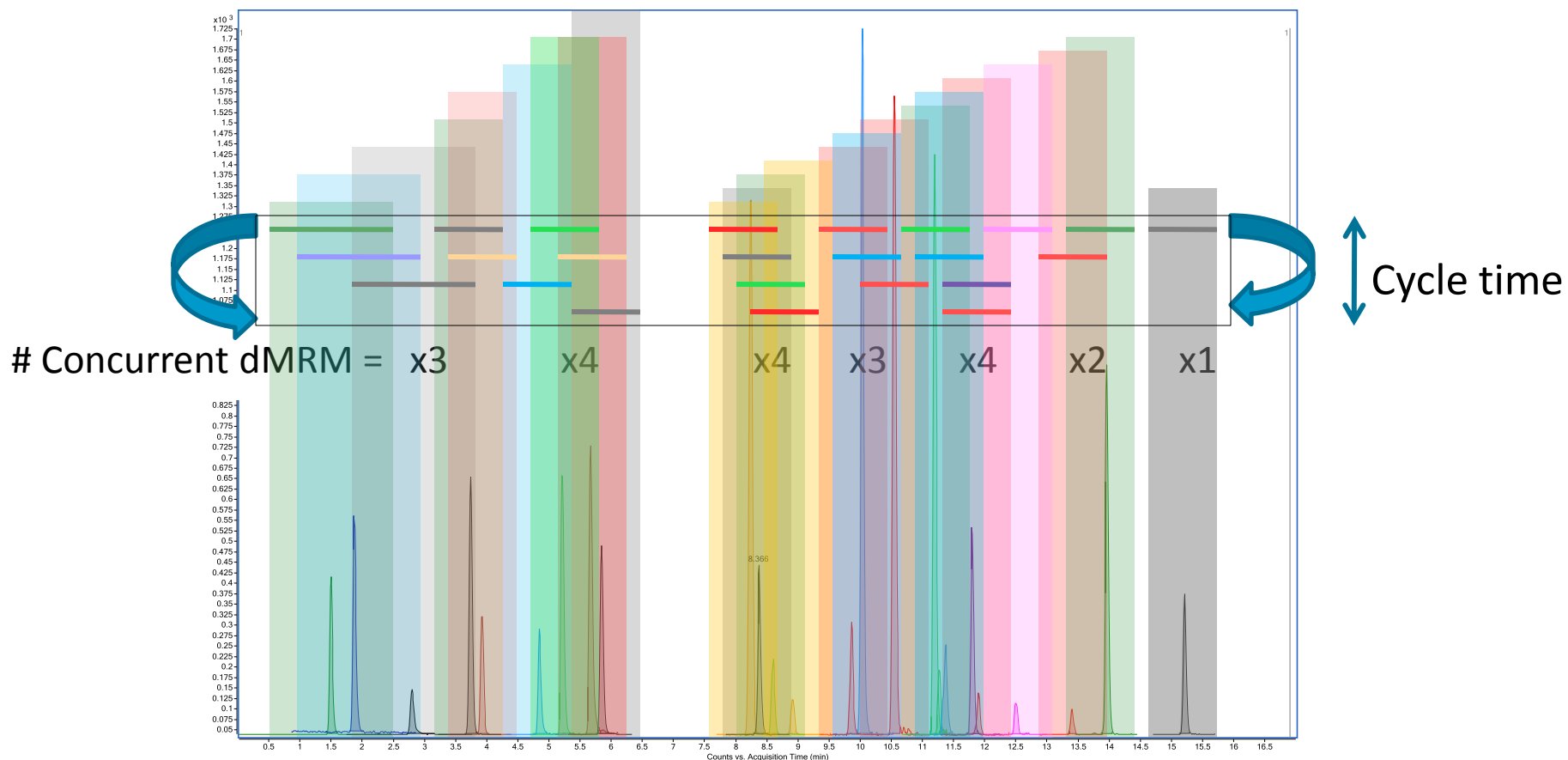
EIC Overlay of 250 pesticide Mix spiked into Tap Water
(500 total transitions, on-column amount 2.5pg) using dynamic MRMs.



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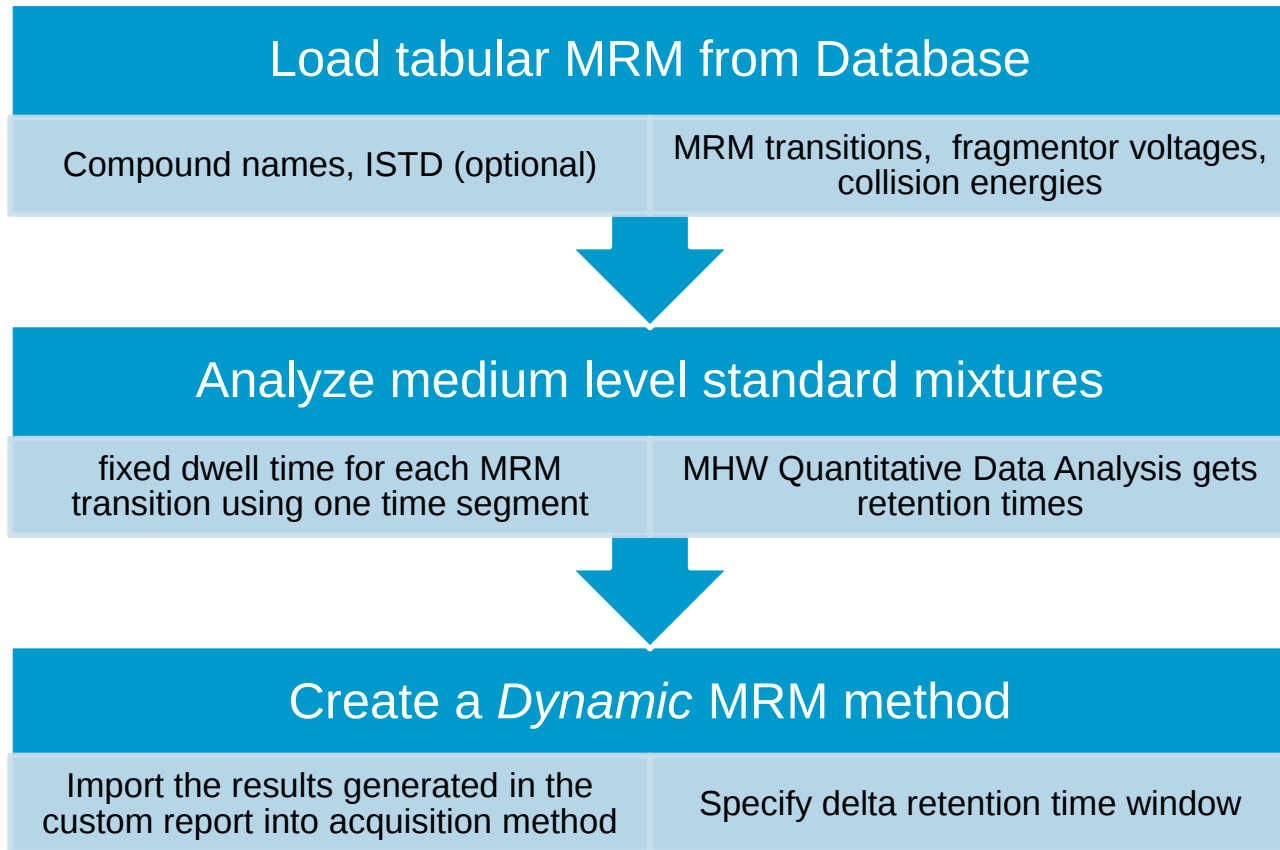


Selection of 24 pesticide transitions from previous slide
To illustrate efficiency of cycle times using dynamic MRMs.



Selection of 24 pesticide transitions from previous slide
To illustrate efficiency of cycle times using dynamic MRMs.

Building a Dynamic MRM Acquisition Method



Agilent 600 Compound Dynamic MRM Database

Database Browser

Database

Filter Compounds

☐ Optimized Compounds

☐ Date From To

☐ Group Name

☐ Project Name

☐ Polarity

Search Compounds

☐ Compound Name

☐ Formula

☐ Method

☐ Show All Records ☐ Show results summary

Compound Information

☐	Compound Name	Group	Formula	Nominal Mass	Vial Number	Project Name
☐	Acetamiprid	Insecticide	C10H11N4Cl	222.07	Vial 1	DefaultProject

☐	Project Name	Method	Polarity	Ion Source	Instrument ID	Date Optimized	Flagged	Flag ID
☐	DefaultProject	E:\MassHunter\Metho	Positive	ESI			☒	

☐	Precursor Ion	Fragmentor	Abundance
☐	223.07	80	

☐	Product Ion	Collision Energy	Abundance
☐	126.01	15	321734
☐	56	15	152499

☐	Compound Name	Group	Formula	Nominal Mass	Vial Number	Project Name
☐	Penconazol	Fungicide	C13H15Cl2N3	283.06	Vial 1	DefaultProject
☐	Quinoxifen	Fungicide	C15H8Cl2FNO	307	Vial 1	DefaultProject
☐	Parathion-methyl	Insecticide	C8H10NO5PS	263	Vial 1	DefaultProject
☐	Indoxacarb	Insecticide	C22H17ClF3N3O	527.07	Vial 1	DefaultProject
☐	Fenfuram	Fungicide	C12H11NO2	201.08	Vial 1	DefaultProject
☐	Dodemorph	Fungicide	C18H35NO	281.27	Vial 1	DefaultProject

Current Database : D:\MassHunter\Databases\MassHunter_Pesticide_DynamicMRM_Database

Refresh Save Import Cancel

Dynamic MRM Acquisition Parameters – *Customized Methods from Database*

The screenshot shows the Agilent software interface for Dynamic MRM Acquisition Parameters. The 'Acquisition' tab is selected, displaying a table of scan segments. A context menu is open over the table, with 'Import from optimizer...' highlighted.

Time segments

#	Start Time	Scan Type	Polarity	Div Valve	Delta EMV	Stored
1	0	Dynamic MRM	Negative	To MS	300	☒

Acquisition | Source | Chromatogram | Instrument | Diagnostics

Scan segments

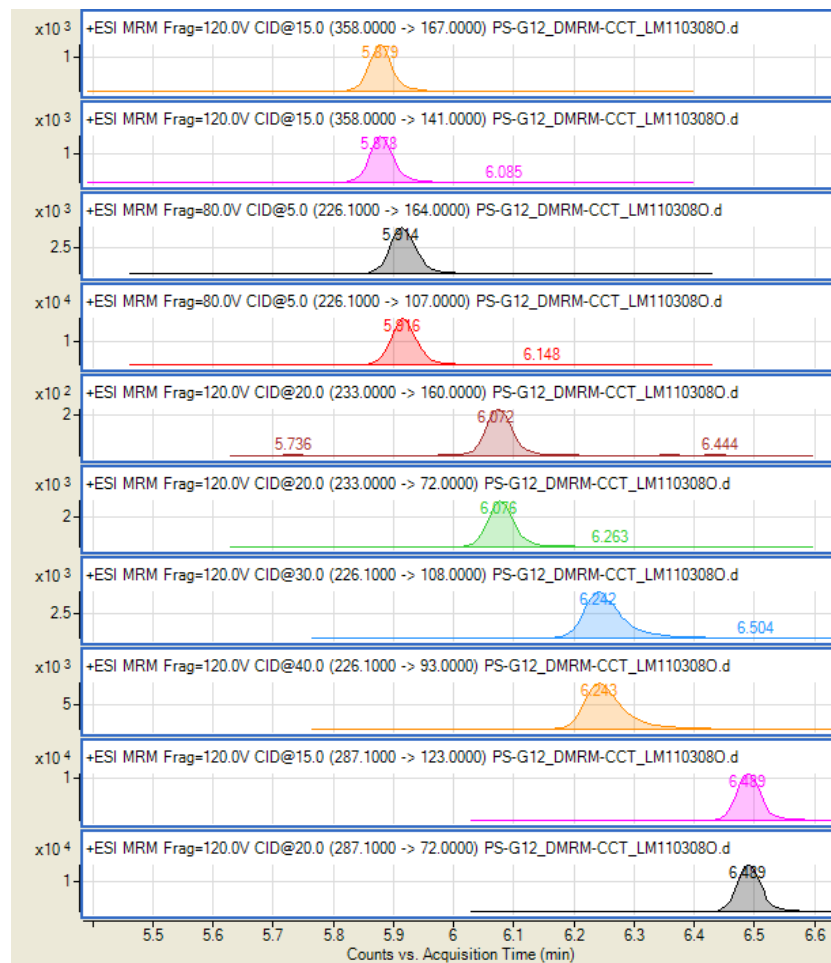
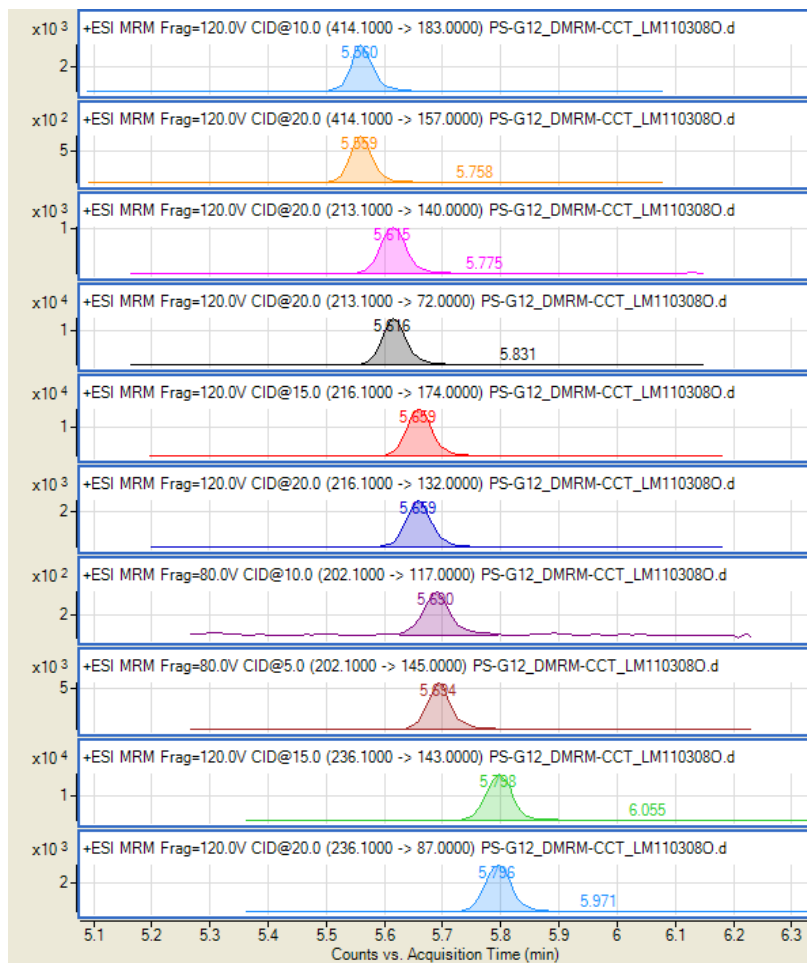
Compound Name	ISTD?	Precursor Ion	MS1 Res	Product Ion	MS2 Res	Fragmentor	Collision Energy	Ret Time (min)	Delta Ret Time
Bentazon	☐	239.1	Unit	132	Unit	80	32	7.117	1
2,4,5-T	☐	252.9	Unit	194.8	Unit	76	9	8.565	1
Silvex	☐	266.9	Unit	194.9	Unit	90	5	9.325	1
Acifluorfen	☐	360	Unit	315.9	Unit	78	5	10.119	1
Dinoseb	☐	239.1	Unit	207	Unit	154	21	11.226	1
Hexaflumuron	☐	459	Unit	438.9	Unit	102	5	11.423	1

Dynamic MRM Parameters

- Add Row
- Delete Row
- Sort
- Import from optimizer...
- Cut
- Copy
- Paste
- Paste from Clipboard
- Fill Down
- Fill Column

Dynamic MRM at the 10 pg level for 100 pesticides

1 min $\Delta R.T.$ - based acquisition window for each MRM transition



Recent Advances in Ultra High Definition Mass Spectrometry with the 1290 Infinity uHPLC System

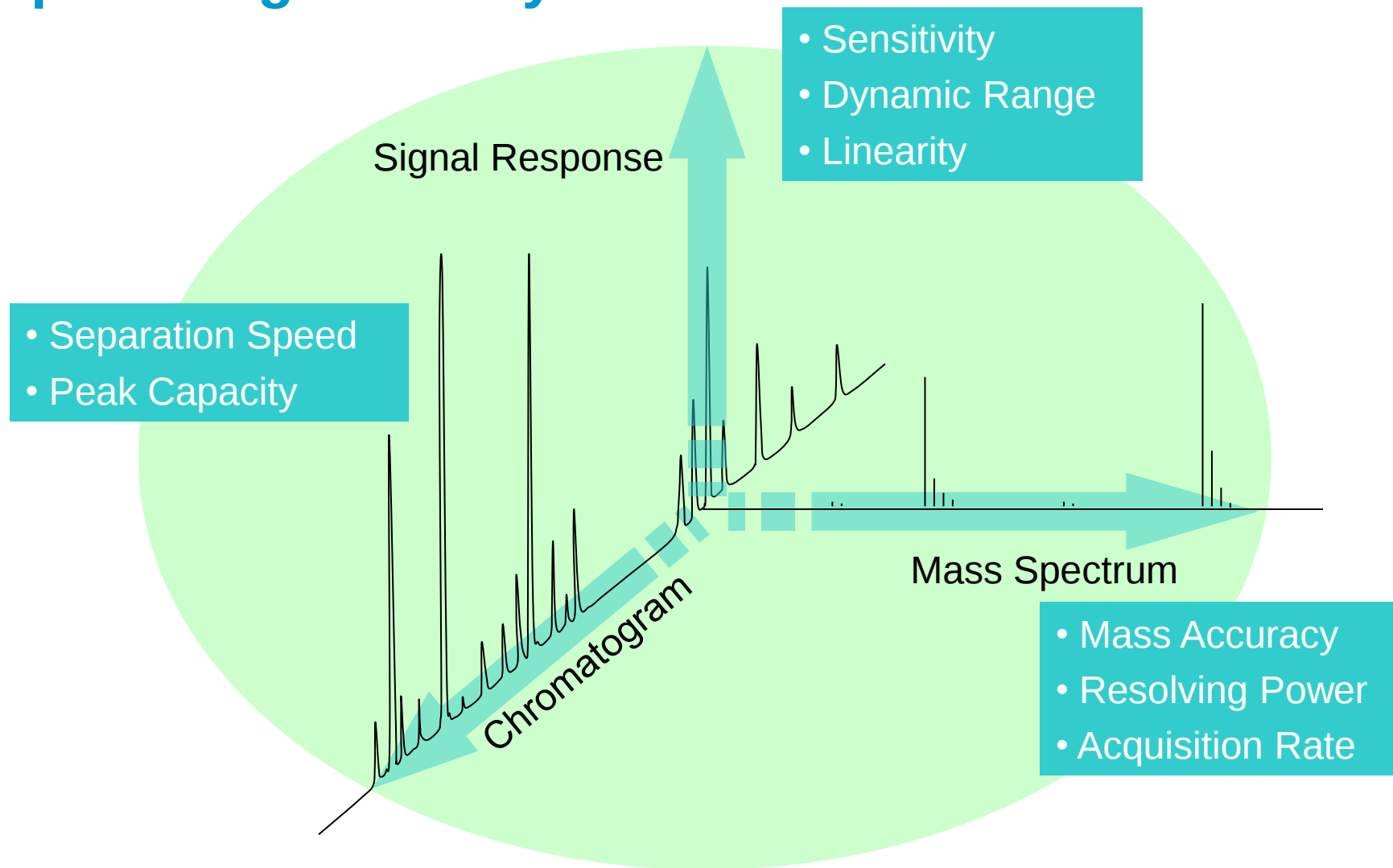
(Agilent 1290 Infinity uHPLC,
Agilent 6230 TOF and 6540 Q-TOF.)

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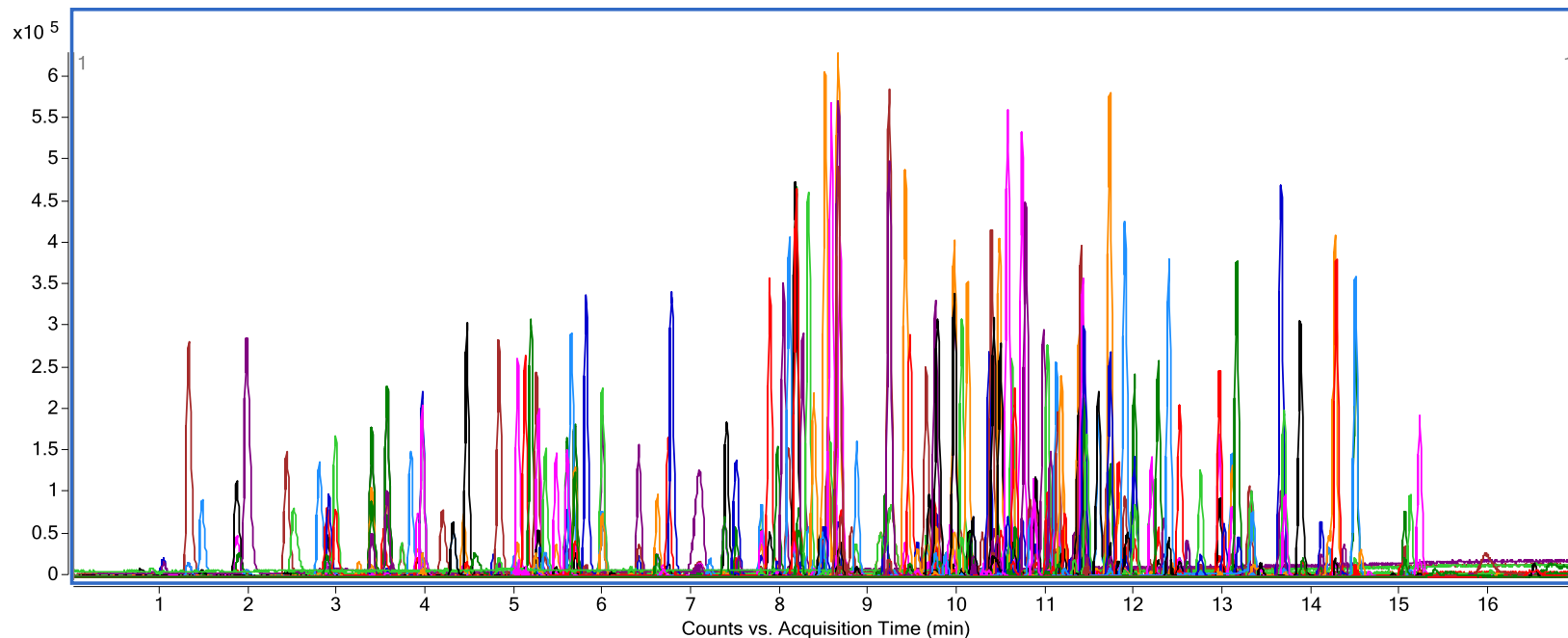


Ultra High Definition

Optimizing all Analytical Dimensions

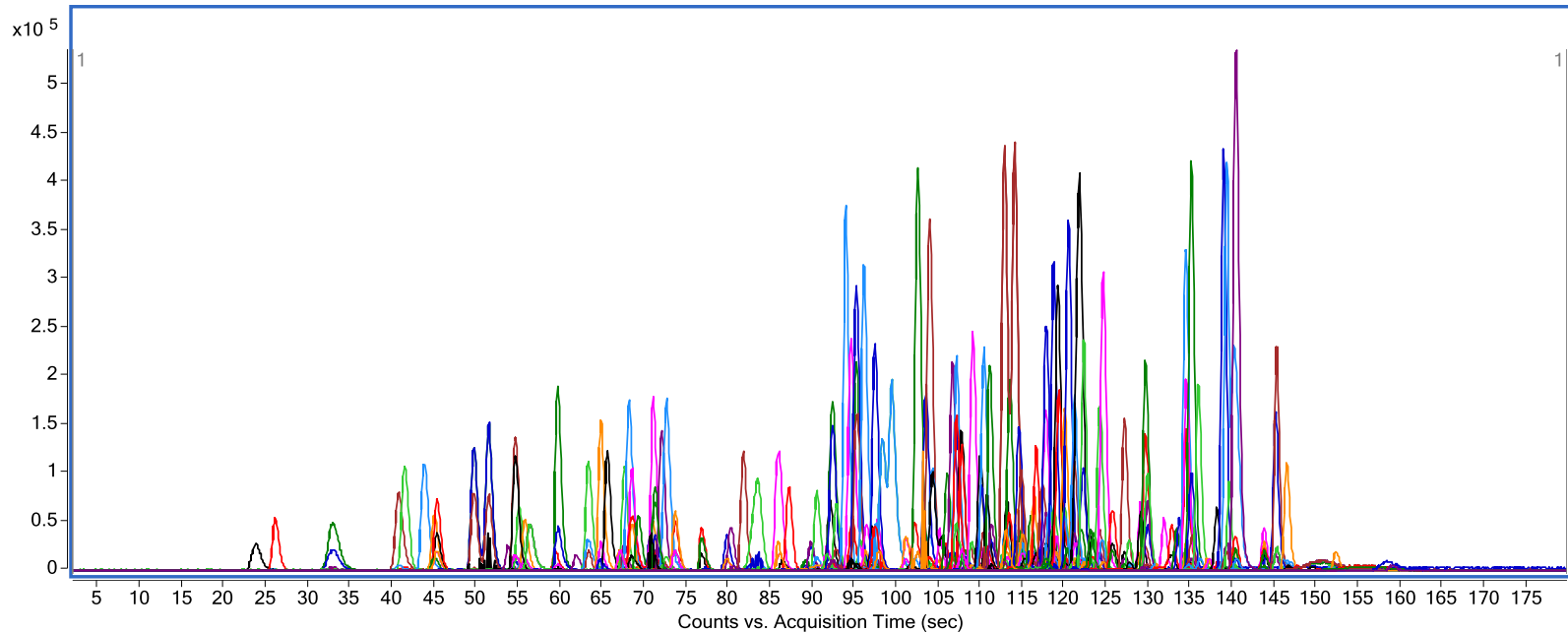


17 minute analysis – 224 Pesticides 6230 TOF/1290 Infinity uHPLC



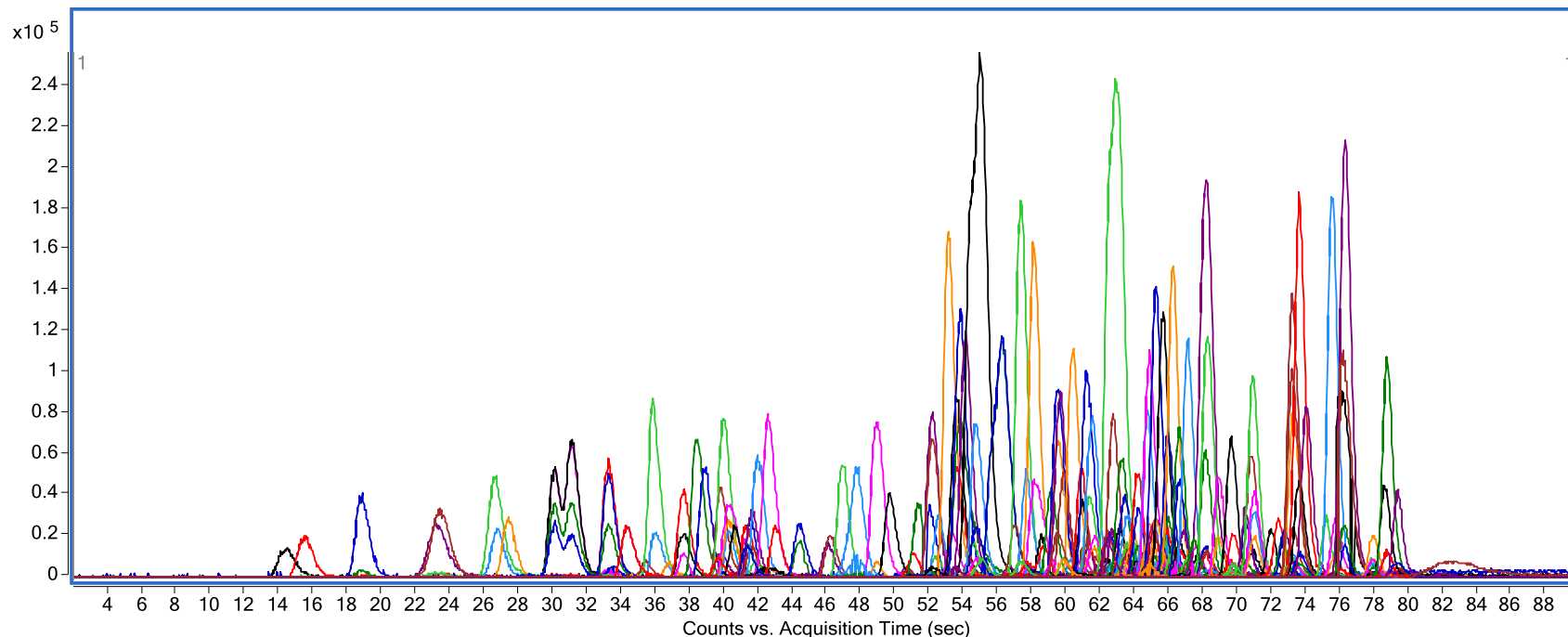
3 Hz Data Acquisition Speed
50pg on-column

3 minute analysis – 224 Pesticides 6230 TOF/1290 Infinity uHPLC



10 Hz Data Acquisition Speed
50pg on-column

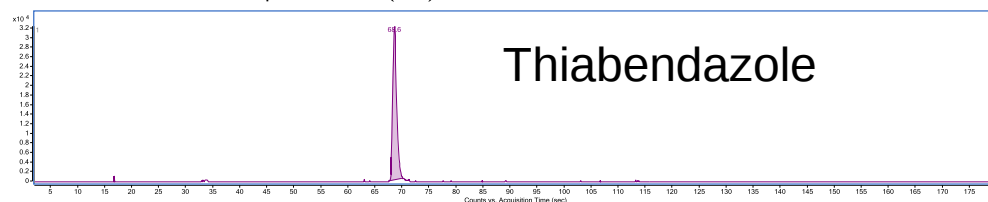
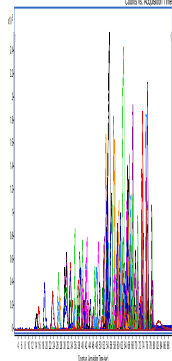
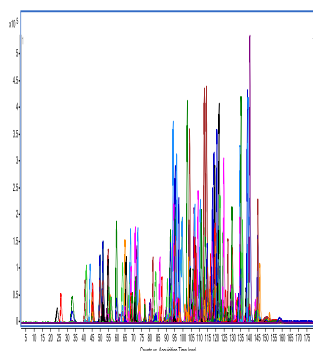
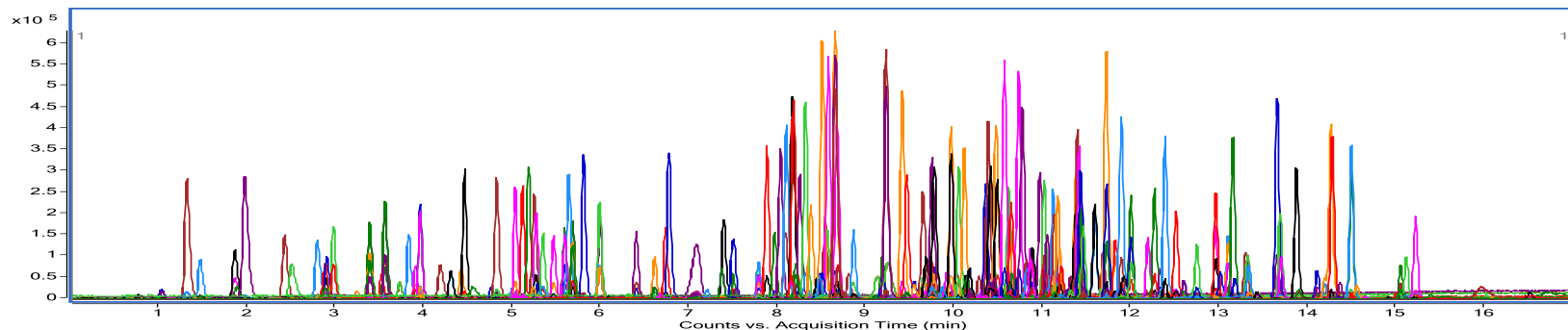
90 second analysis – 224 Pesticides 6230 TOF/1290 Infinity uHPLC



20 Hz Data Acquisition Speed
50pg on-column

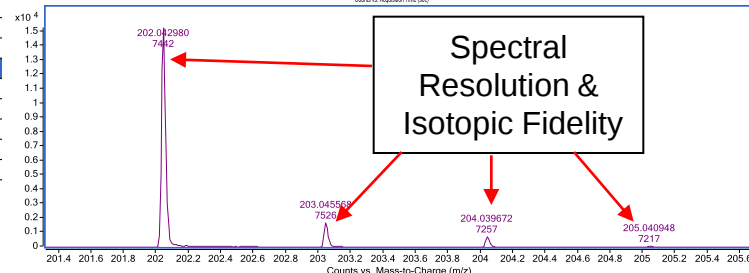
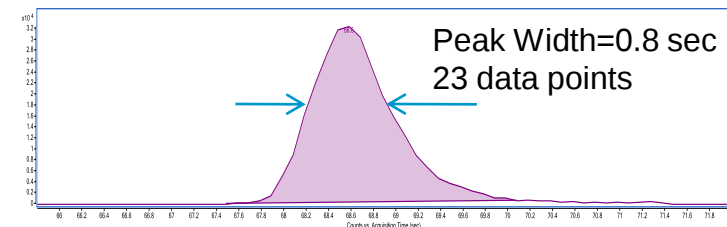
Agilent 6230 TOF Pesticide Screening

224 pesticides, 50 pg on column



Compound List					
Cpd	Name	RT	Mass	DW (Tgt. ppm)	DW (Tgt. mDa)
191	Tebufenpyrad	123	333.1613	1.43	0.48
174	Tebupirimfos	138	318.1172	1.65	0.52
145	Tebupirimfos oxone	107.1	302.1397	0.36	0.11
77	Tebuthiuron	81.7	228.1044	-0.21	-0.05
115	Terbufos	120	288.0437	-1.59	-0.46
74	Terbufos	97.7	225.1589	-0.05	-0.01
28	Terbutylazine-desethyl	54.7	201.0788	3.46	0.7
91	Terbutynl	112.8	241.1362	0.5	0.12
215	Tetraconazol	110.8	371.0216	0.24	0.09
25	Thiabendazole	68.6	201.0357	-1.81	-0.36
94	Thiabendazole	77	252.0233	-1.38	-0.35
118				0.08	
127				-11.66	
167				-0.16	
169	Tribufos	146.3	314.0964	0.76	0.24
161	Triurilazole	77	189.0794	-1.97	-1.18

-1.8 ppm



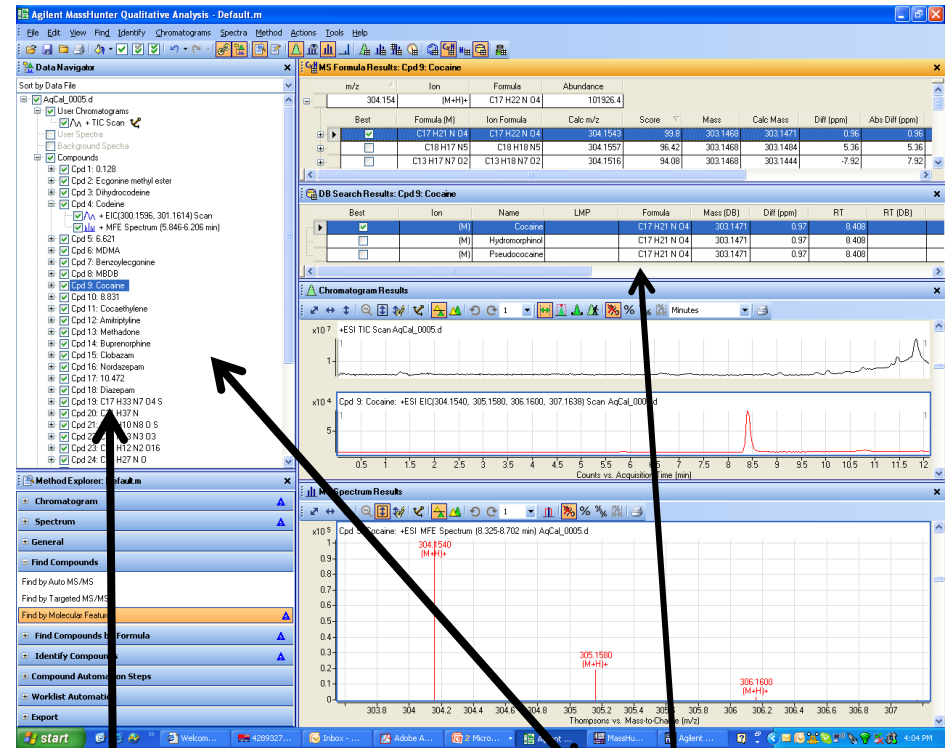
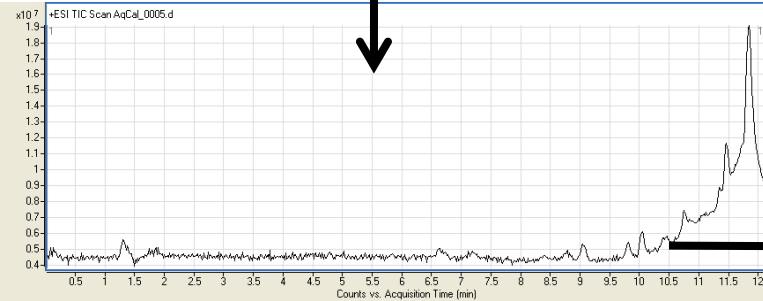
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Agilent TOF Screening Solution (PCD)



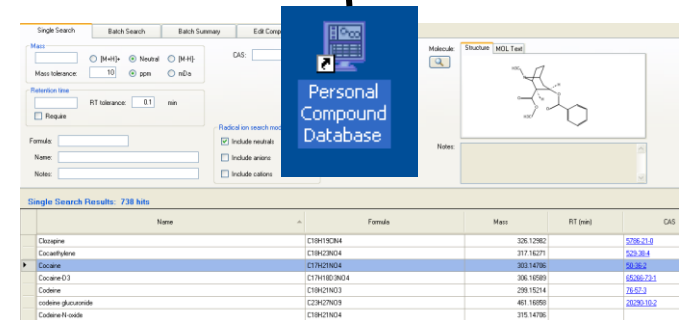
Agilent 6230 TOF

- 20K resolution
- 5 decades in-scan
- Low Pg Sensitivity



Molecular Feature
Extractor

Find by
Formula



Agilent Technologies

PCD Search Results For 50pg Forensic Spiked Blood

– PCD Search Confirmed compound 9 as Cocaine

Data Navigator

Sort by Data File

- [-] AqCal_0005.d
 - [-] User Chromatograms
 - [-] User Spectra
 - [-] Background Spectra
 - [+] Compounds
 - [x] Cpd 1: 0.128
 - [x] Cpd 2: Ecgonine methyl ester
 - [x] Cpd 3: Dihydrocodeine
 - [x] Cpd 4: Codeine
 - [x] Cpd 5: 6.621
 - [x] Cpd 6: MDMA
 - [x] Cpd 7: Benzylecgonine
 - [x] Cpd 8: MBDB
 - [x] **Cpd 9: Cocaine**
 - [x] Cpd 10: 8.831
 - [x] Cpd 11: Cocaethylene
 - [x] Cpd 12: Amitriptyline
 - [x] Cpd 13: Methadone
 - [x] Cpd 14: Buprenorphine
 - [x] Cpd 15: Clobazam
 - [x] Cpd 16: Nordazepam
 - [x] Cpd 17: 10.472
 - [x] Cpd 18: Diazepam
 - [x] Cpd 19: C17 H33 N7 O4 S
 - [x] Cpd 20: C21 H37 N
 - [x] Cpd 21: C13 H10 N8 O6
 - [x] Cpd 22: C20 H13 N3 O3
 - [x] Cpd 23: C18 H13 N3 O16

Compound List

Cpd	Name	RT	Mass	Mass (MFG)	Diff (MFG, ppm)	Algorithm	Score (DB)	Formula (MFG)	Score (MFG)	Hits (DB)	Ions
1		0.128	120.1787			Find by Molecular Feature				2	
2	Ecgonine methyl est...	1.404	199.1208			Find by Molecular Feature	99.88			1	4
3	Dihydrocodeine	5.849	301.1677			Find by Molecular Feature	99.67			2	4
4	Codeine	6.975	299.1523	299.1521	-0.68	Find by Molecular Feature	83.51	C18 H21 N O3	84.09	2	2
5		6.621	149.1192			Find by Molecular Feature				3	
6	MDMA	6.727	193.1101			Find by Molecular Feature	99.72			3	3
7	Benzylecgonine	7.141	289.131			Find by Molecular Feature	99.55			1	4
8	MBDB	7.211	207.1256			Find by Molecular Feature	87.15			2	2
9	Cocaine	8.408	303.1468	303.1471	0.97	Find by Molecular Feature	99.68	C17 H21 N O4	99.8	3	4
10		8.431	162.1253			Find by Molecular Feature				3	
11	Cocaethylene	9.054	317.1625			Find by Molecular Feature	98.12			1	4
12	Amitriptyline	9.312	277.183			Find by Molecular Feature	72.88			3	2
13	Methadone	10.045	309.2093			Find by Molecular Feature	98.44			1	3
14	Buprenorphine	10.053	467.3033	467.3036	0.6	Find by Molecular Feature	99.51	C29 H41 N O4	99.54	1	4
15	Clobazam	10.397	300.0662			Find by Molecular Feature	98.97			2	5
16	Nordazepam	10.41	270.0556			Find by Molecular Feature	99.53			1	4
17		10.472	213.0821			Find by Molecular Feature				4	
18	Diazepam	10.759	284.0718			Find by Molecular Feature	96.95			1	5

DB Search Results: Cpd 9: Cocaine

Best	Ion	Name	Formula	Mass (DB)	Diff (ppm)	RT	RT (DB)	RT Diff	Abund Match	Spacing Match	Mass Match	RT Match
[x]	(M)	Cocaine	C17 H21 N O4	303.1471	0.97	8.408			99.94	98.99	99.87	
[]	(M)	Hydromorphenol	C17 H21 N O4	303.1471	0.97	8.408			99.94	98.99	99.87	
[]	(M)	Pseudococaine	C17 H21 N O4	303.1471	0.97	8.408			99.94	98.99	99.87	

Focused AMRT Databases

(Personal Compound Databases (PCD))

Pesticides, Part # (G6854AA):

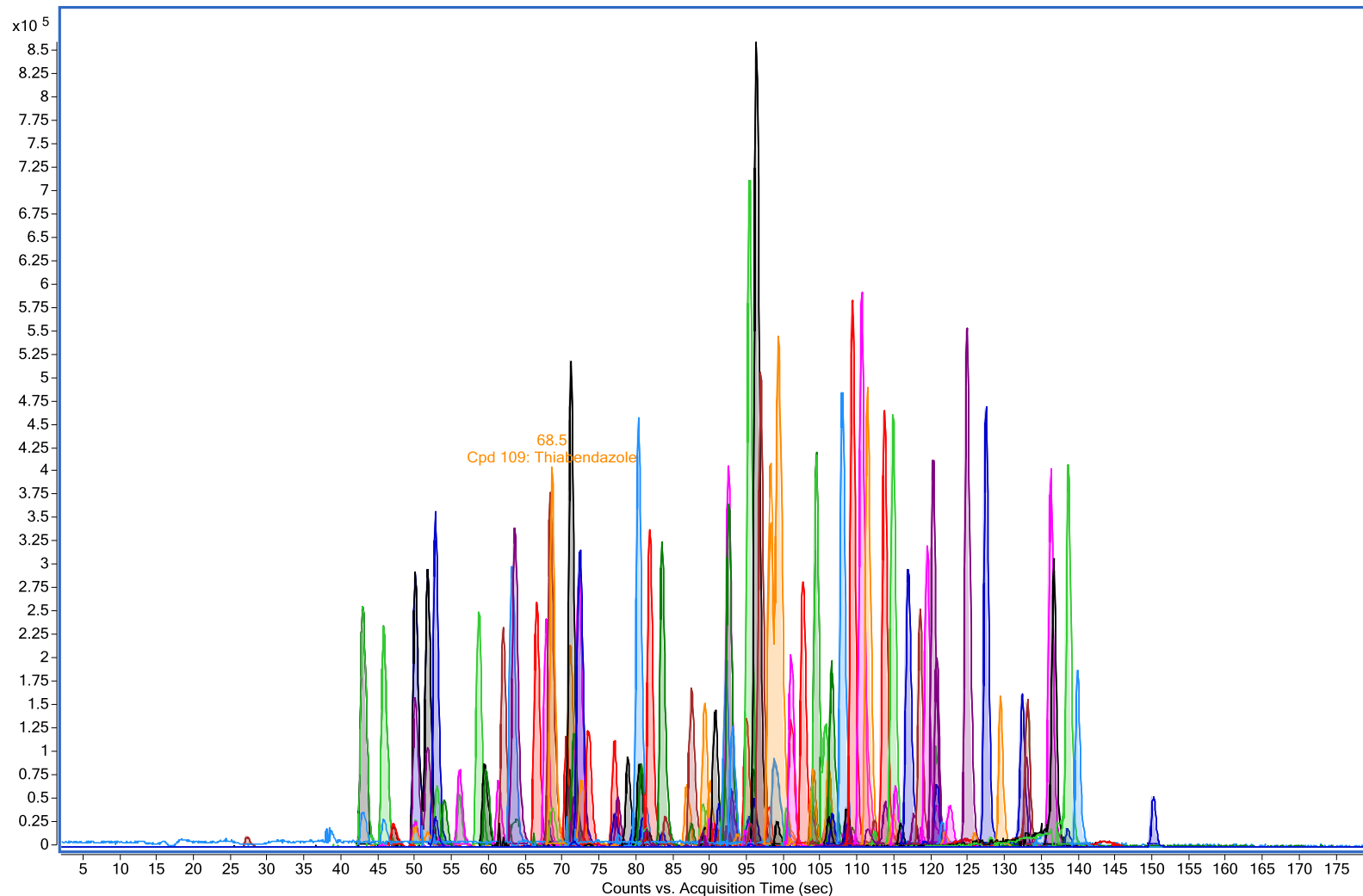
- 1600 analyte content (formulae & accurate mass)
- Structures
- CAS Numbers & Links to external website (pubchem)

Forensic & Toxicology, Part # (G6855AA):

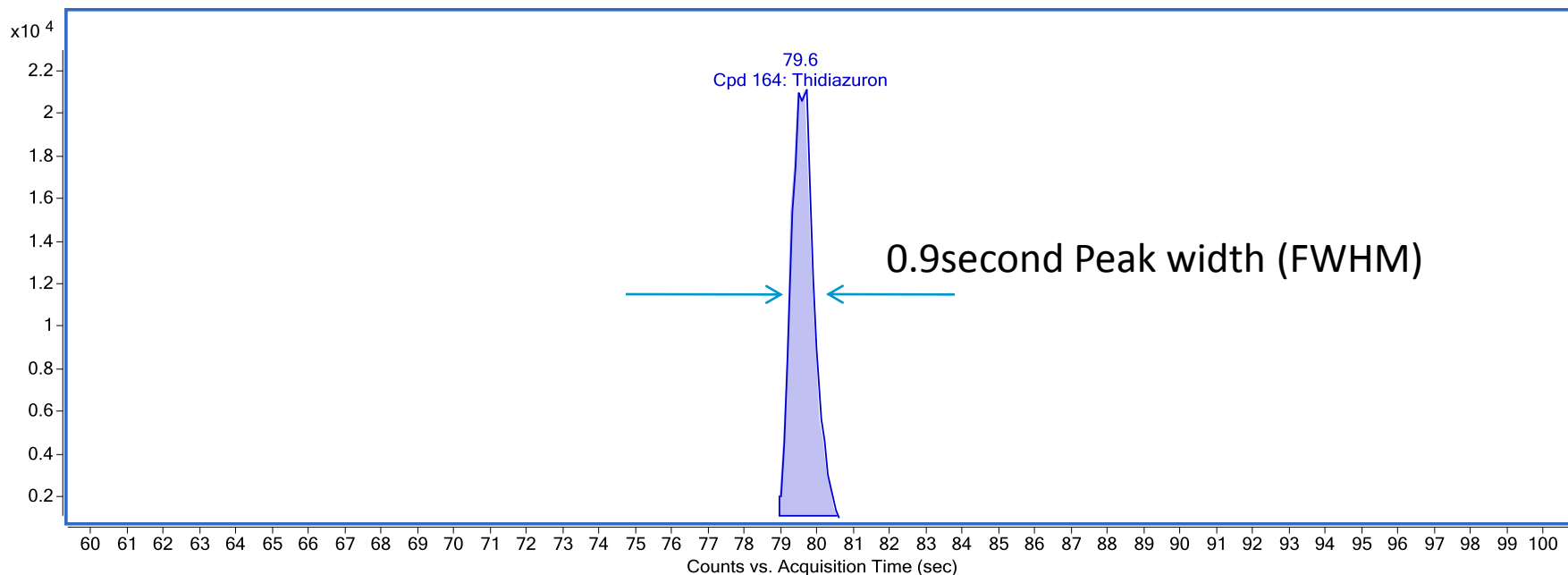
- ~7000 analytes content (formulae & accurate mass)
- Structures
- CAS Numbers & Links to external website (pubchem)



PESTICIDE Non-target Screen, 6230 TOF/1290 uHPLC
3 minute analysis, Zorbax Eclipse + C18 2.1x100mm, 1.8um
100pg on-column, 10 Hz Acquisition.



**3 minute analysis, Zorbax Eclipse + C18 2.1x100mm, 1.8um
6230 TOF/1290uHPLC, 100pg on-column, 10 Hz Acquisition.**



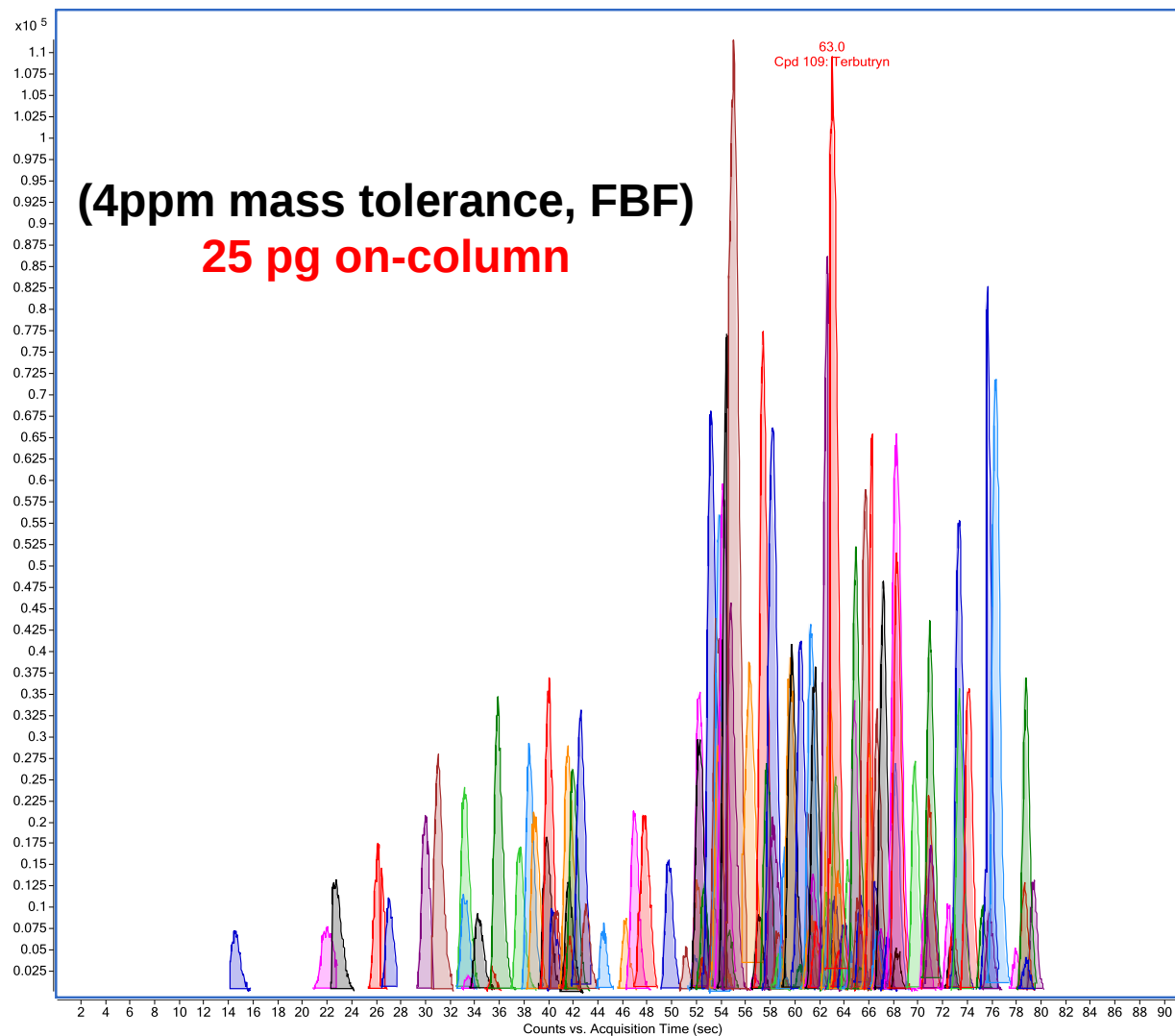
1290 Infinity/6230 TOF

Find By Formula PCD Search

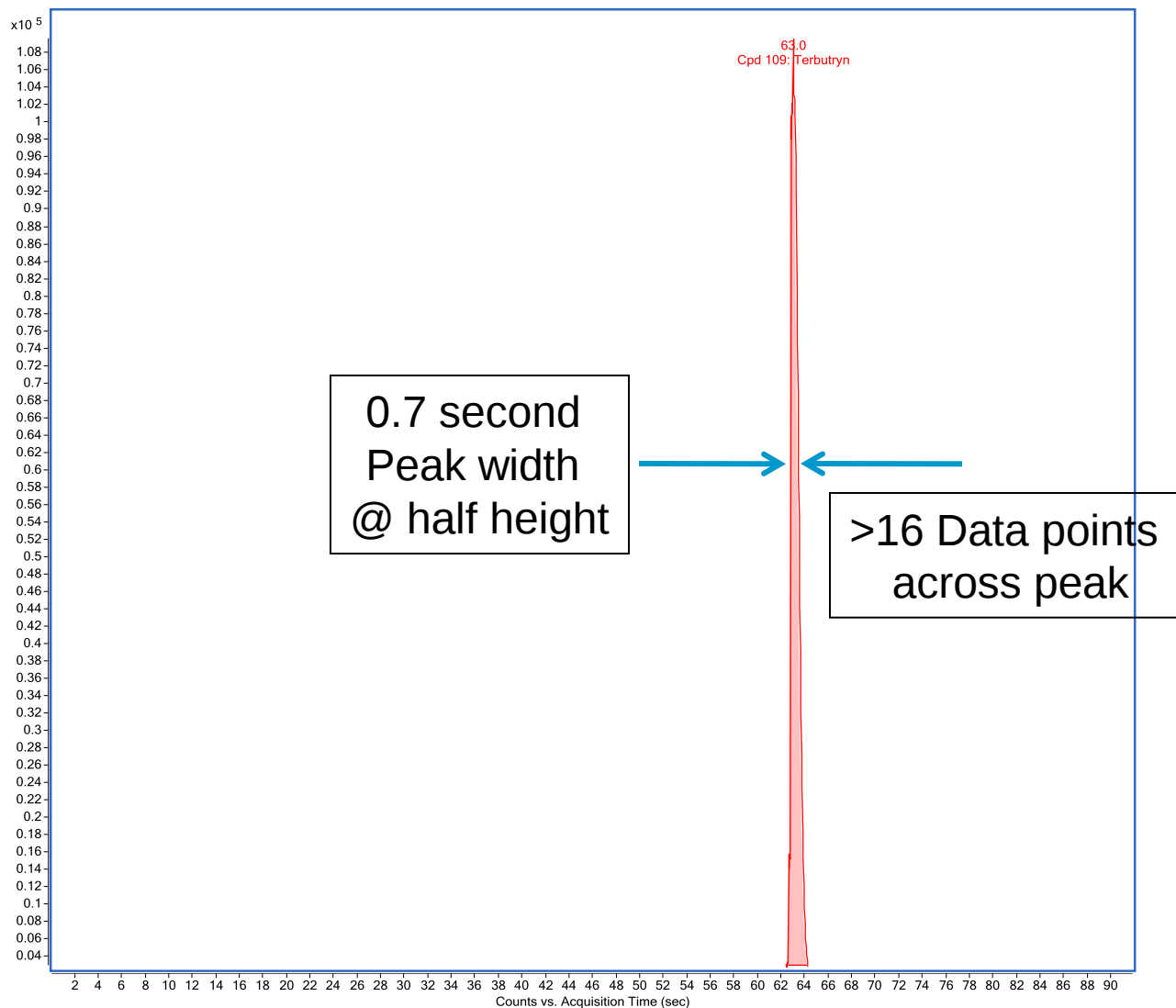
3-minute 1290/6230 TOF Screening (224 +ve Pesticide suite)		
Analyte amount (on-column)	# of compounds identified (%)	Average Mass Accuracy (ppm)
500fg	45 (20%)	0.24
1.25pg	96 (43%)	0.51
2.5pg	124 (55%)	0.49
5pg	163 (73%)	0.12
25pg	202 (90%)	0.70
50pg	217 (97%)	0.06
125pg	224 (100%)	0.60



Non-target Screen (6230 TOF/1290 uHPLC), 202/224 (90%) Pesticides Identified 90 second Analysis time 20Hz (2GHz mode)



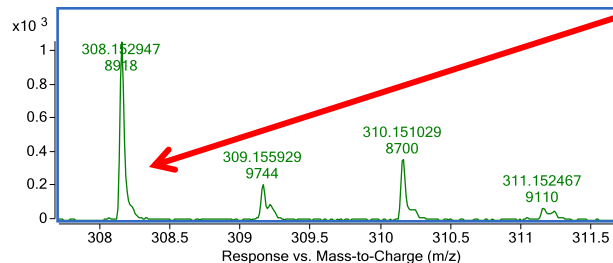
Terbutryn 90 second Analysis time 20Hz (2GHz mode) Zorbax Eclipse Plus C18 (2.1x50mm x 1.8um) 6230TOF/1290 UHPLC



6230 TOF Speed, Mass Resolution & Dynamic Range – Example: Tebuconazole (2GHz mode)

Tebuconazole (on-column amount)	3 min analysis 10Hz acquisition		1.5 min analysis 20Hz acquisition	
	Mass Error (ppm)	Mass Resolution	Mass Error (ppm)	Mass Resolution
500pg on-column	-0.9	8542	1.46	8546
250pg on-column	-1.79	8626	-1.02	8617
125pg on-column	-0.8	8711	1.82	8803
50pg on-column	-1.77	8822	-0.06	8922
25pg on-column	-2.03	8749	1.51	8918
Peak Width FWHM (sec)	0.8		0.5	

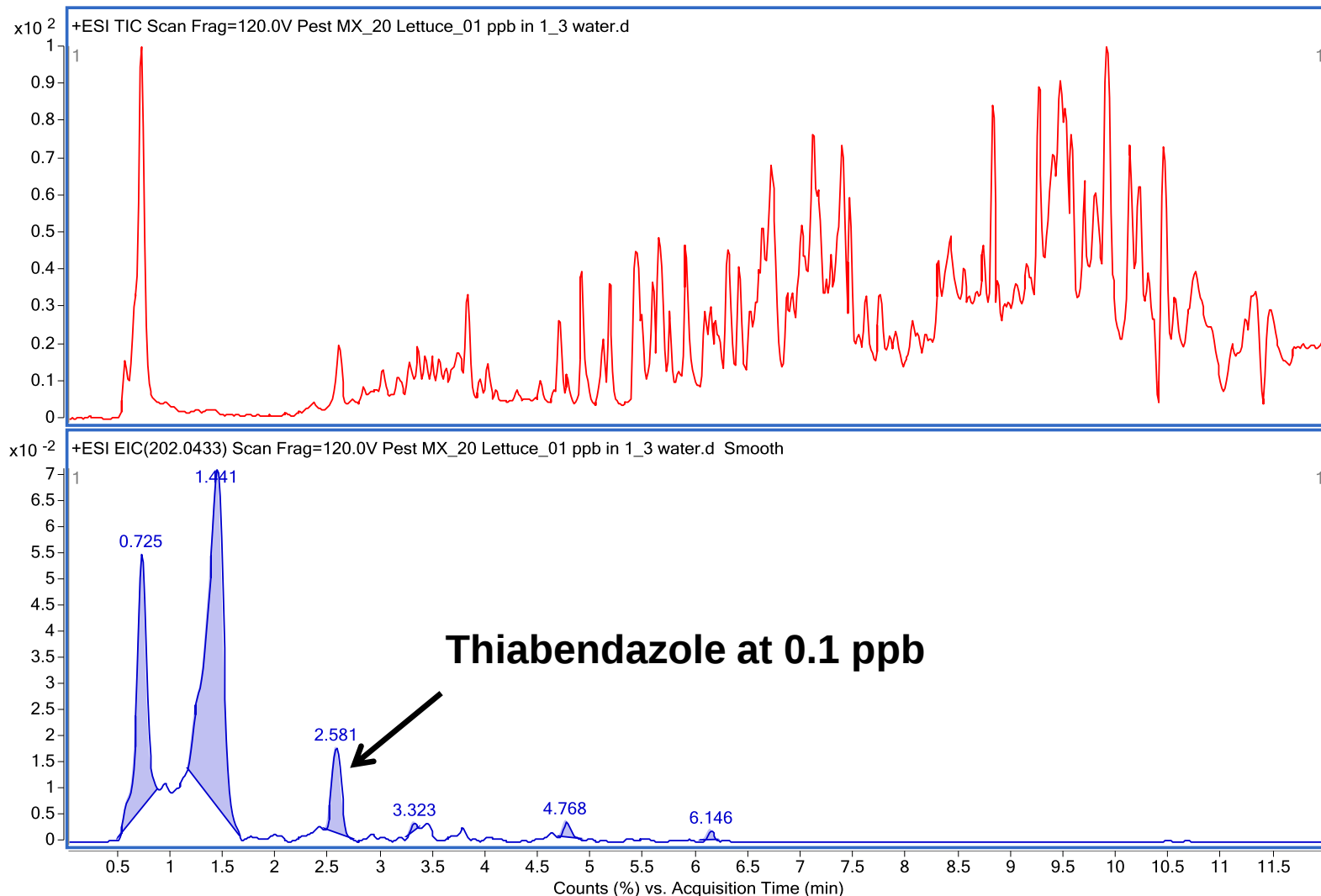
$[M+H]^+ = 308.152947$



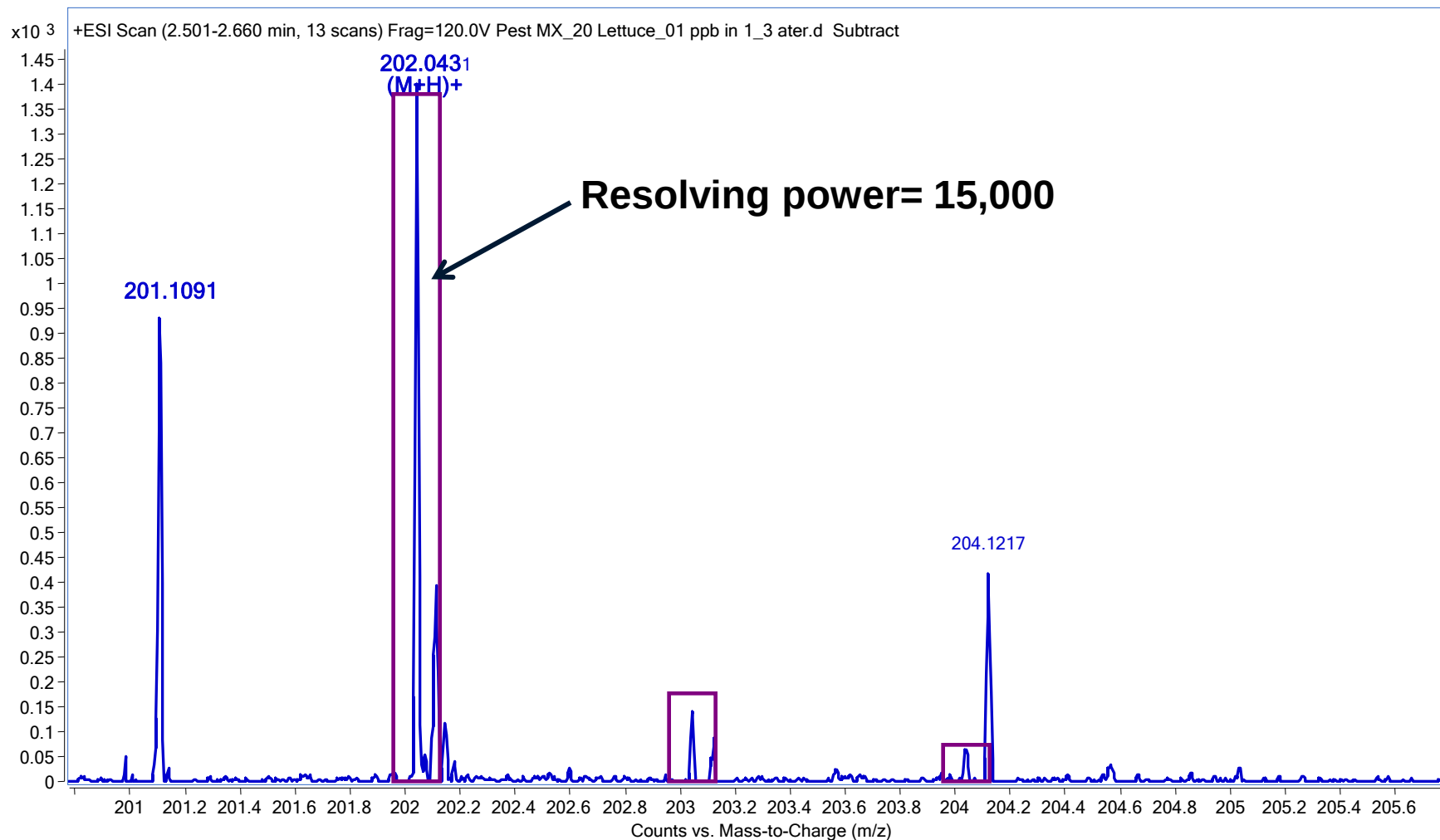
6230 TOF Speed, Mass Resolution & Dynamic Range – Example: Tebuconazole (4GHz mode)

Tebuconazole (on-column amount)	3 min analysis 10Hz acquisition		1.5 min analysis 3Hz acquisition	
	Mass Error (ppm)	Mass Resolution	Mass Error (ppm)	Mass Resolution
250pg on-column	-0.26	17765	-0.2	17949
125pg on-column	0.4	17314	0.1	17432
50pg on-column	-0.31	16959	0.2	17226
25pg on-column	-0.35	17122	0.03	17119
5pg on-column	-1.42	16177	0.03	16499
2.5pg on-column	ND	ND	-1.13	16262
Peak Width FWHM (sec)	0.8		0.5	

Extracted Ion Chromatogram m/z 202.0433±0.002 of Lettuce Extract



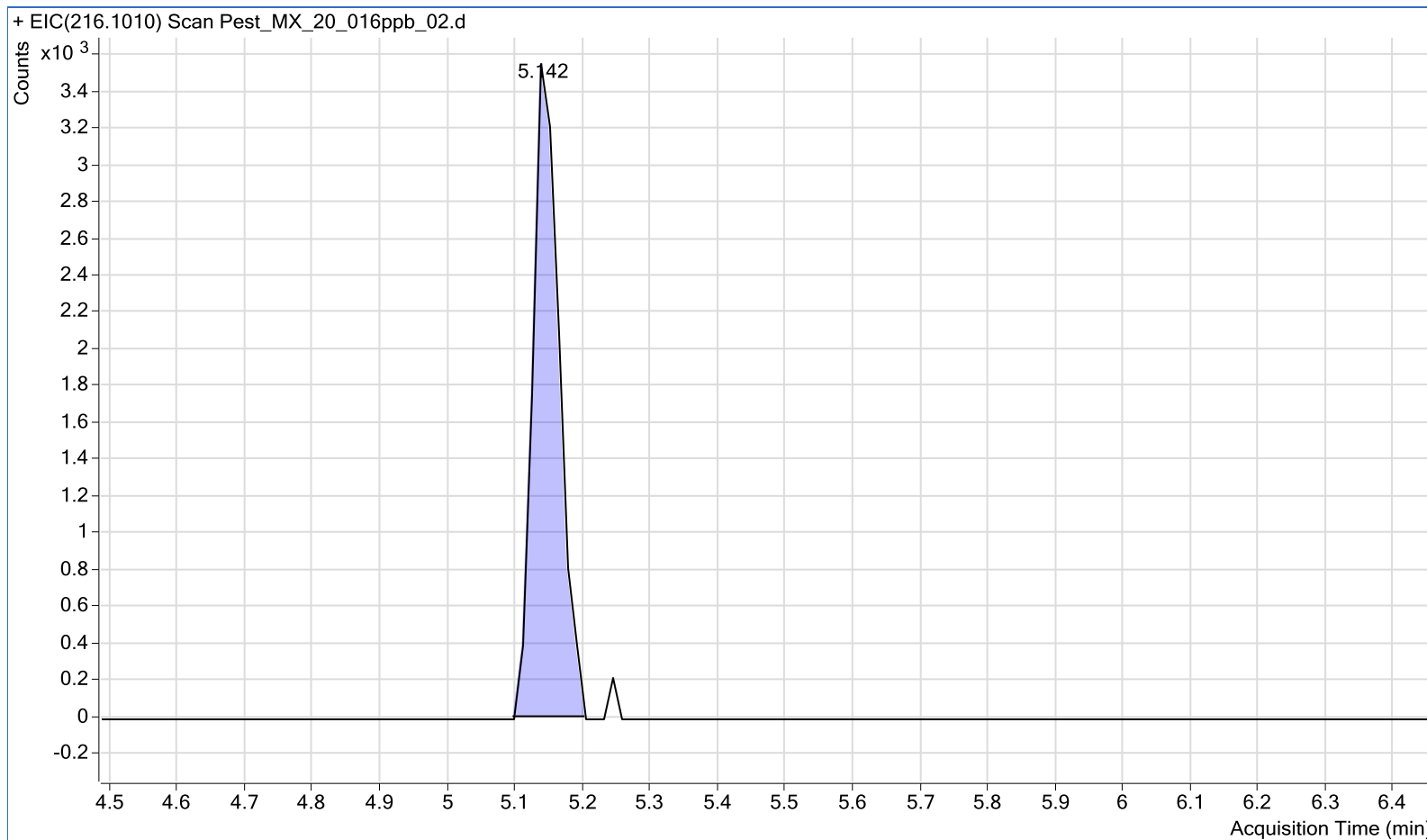
Spectrum of Thiabendazole at 0.1 ppb in Extracted Pepper



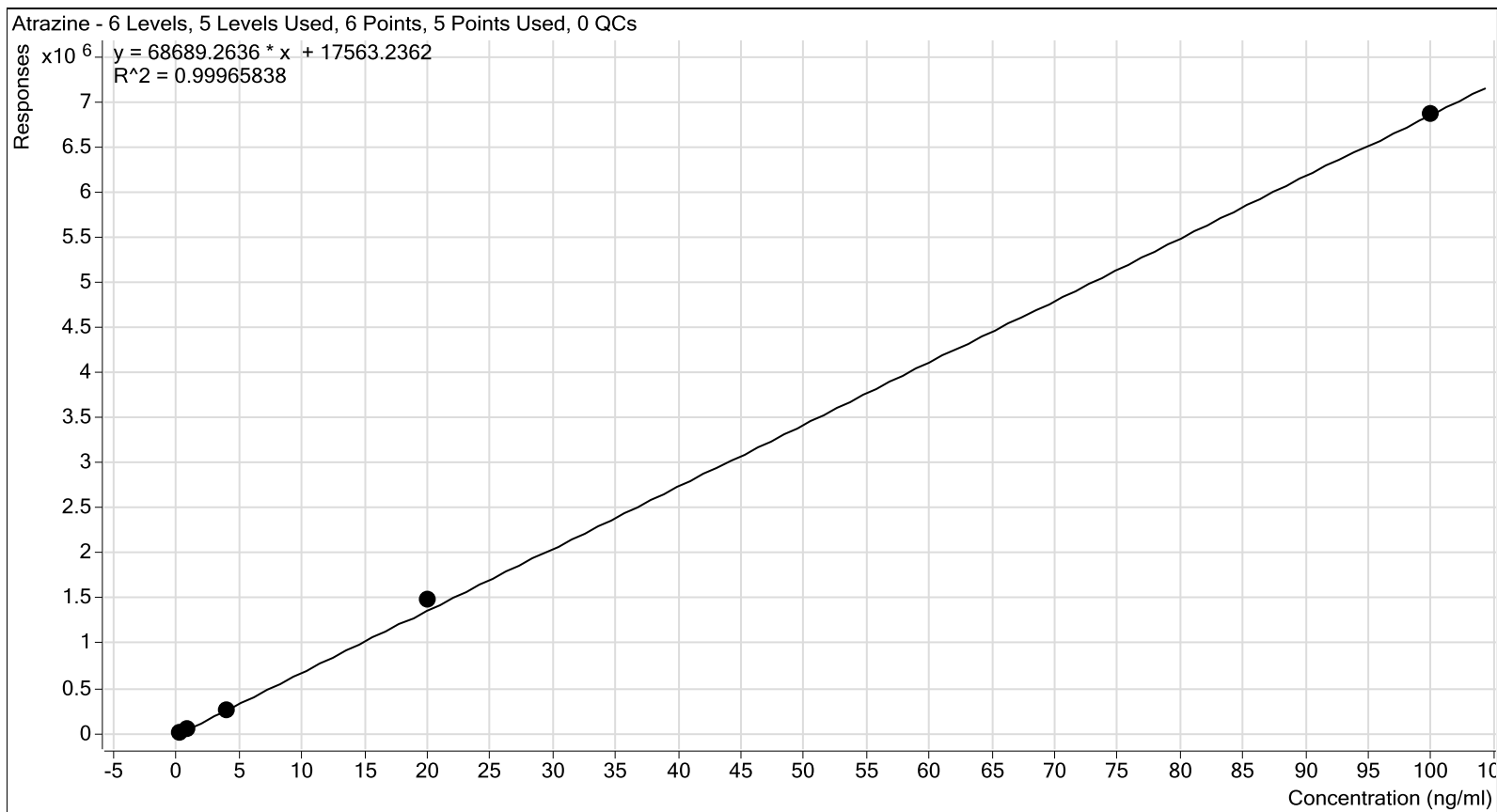
Calculated Formula of Thiabendazole at 0.1 ppb in Extracted Pepper

MS Formula Results: + Scan (2.501-2.660 min) Sub										
m/z	Ion	Formula	Abundance							
202.0431	(M+H)+	C10 H8 N3 S	1415.4							
Best	Formula (M)	Ion Formula	Score	Cross Score	Calc m/z	Diff (ppm)	Mass Match	Abund Match	Spacing Match	
☒	C10 H7 N3 S	C10 H8 N3 S	95.86		202.0433	1.3	99.31	94.72	90.32	
Isotope	Abund Sum%	Calc Abund Sum%	m/z	Calc m/z	Diff (ppm)					
1	86.84	84.73	202.0431	202.0433	1.28					
2	9.16	10.84	203.0447	203.0459	5.87					
3	4.01	4.43	204.0382	204.0405	11.11					
Best	Formula (M)	Ion Formula	Score	Cross Score	Calc m/z	Diff (ppm)	Mass Match	Abund Match	Spacing Match	
☐	C6 H7 N3 O5	C6 H8 N3 O5	59.68		202.0458	13.75	45.89	81.12	61.54	
☐	C12 H8 Cl N	C12 H9 Cl N	59.62		202.0418	-6.36	84.62	0.31	80.79	
☐	C4 H12 Cl N3 O...	C4 H13 Cl N3 O2 S	55.62		202.0412	-9.62	68.31	0.02	96.94	
☐	C9 H12 Cl N S	C9 H13 Cl N S	51.04		202.0452	10.4	64.03	0.05	86.24	
☐	C13 H3 N3	C13 H4 N3	48.82		202.04	-15.46	37.32	64.34	53.17	

Atrazine at 0.16 ppb (0.8 pg on-column)



Atrazine 0.16 ppb to 100



Agilent's New 6540 Ultra High Definition Accurate Mass Q-TOF

Exceptional accurate mass, sensitivity, dynamic range and resolution ... perfect match for 1290 Infinity UHPLC

500 ppb mass accuracy

femtogram sensitivity

5 decades dynamic range

40,000 resolving power

20 Spectra/sec

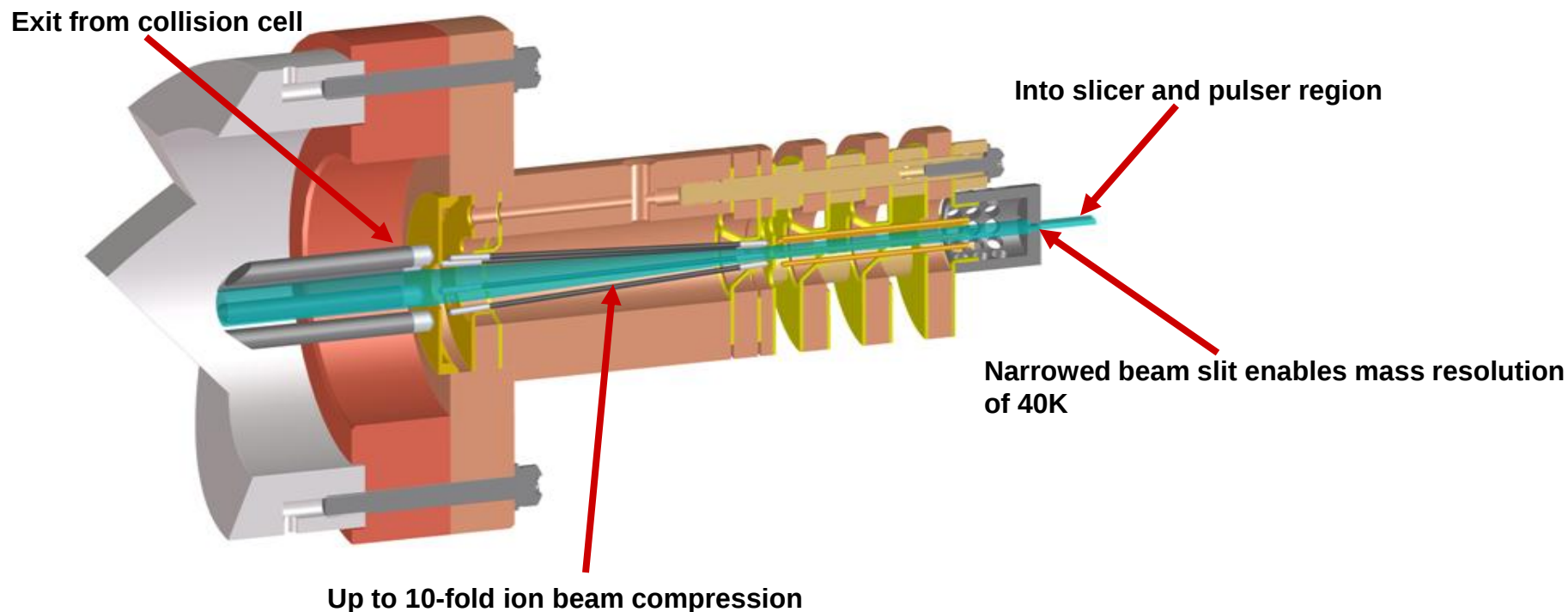
Excellent Linearity and Isotopic Fidelity

Supports Agilent Jet Stream and HPLC-Chip



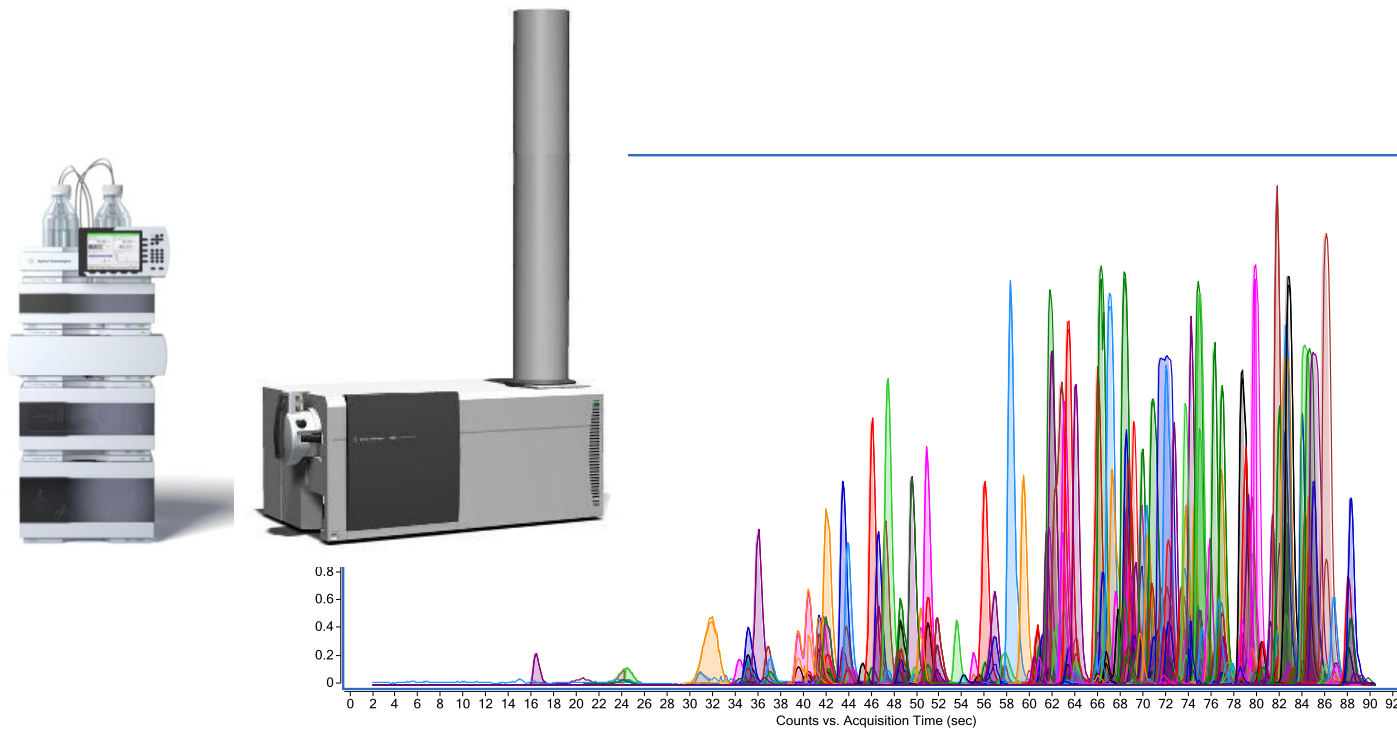
Ion Beam Compression (IBC)* Technology Drives Higher Resolution

Compressed and cooled ion beam ensures the best sensitivity performance in high resolution mode

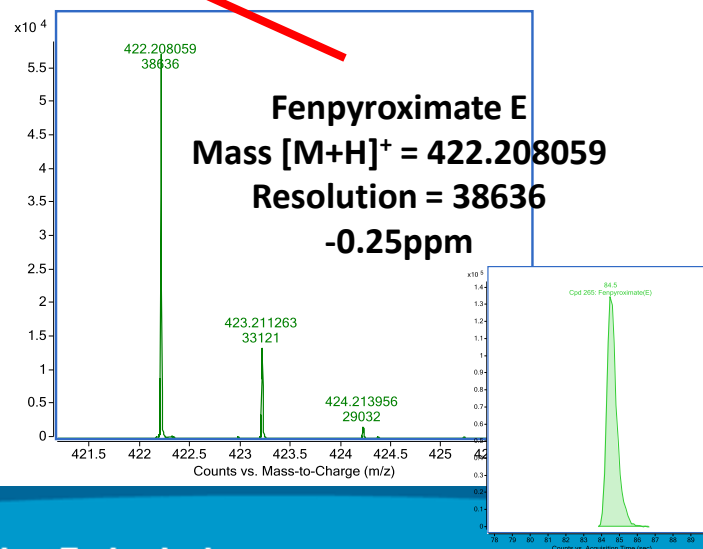
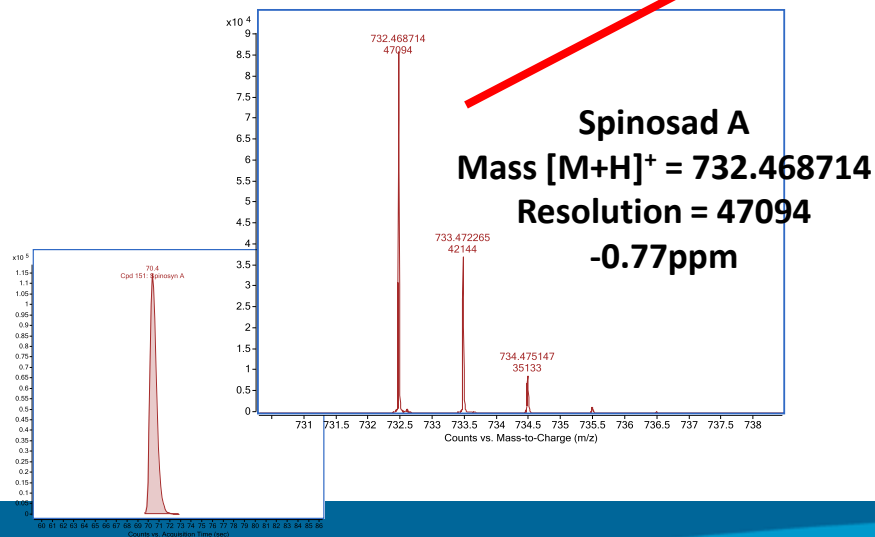
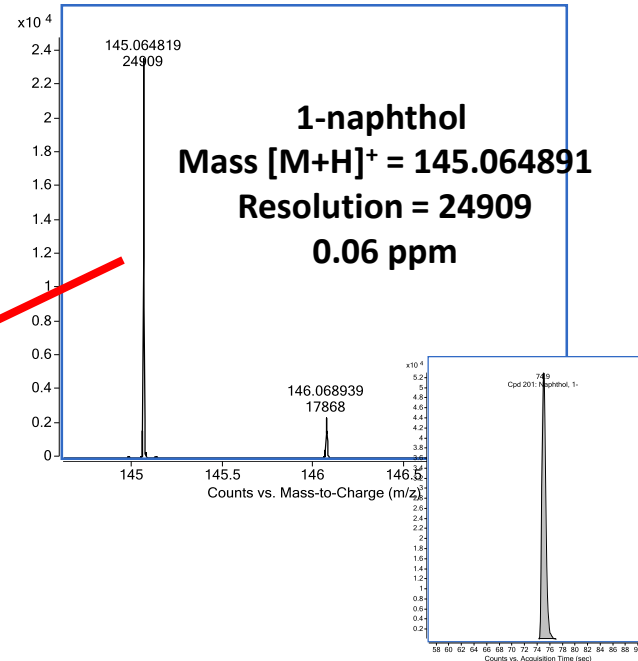
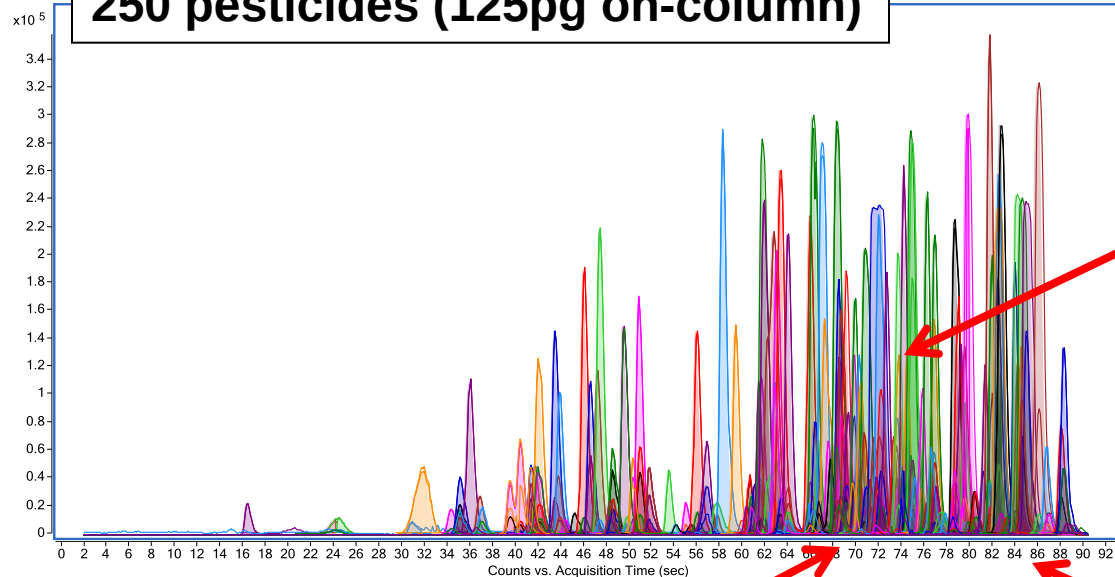


* Patent pending

Pesticide Screen 1290 Infinity/6540 Q-TOF



1290/6540 **90sec** analysis 250 pesticides (125pg on-column)



Agilent Technologies

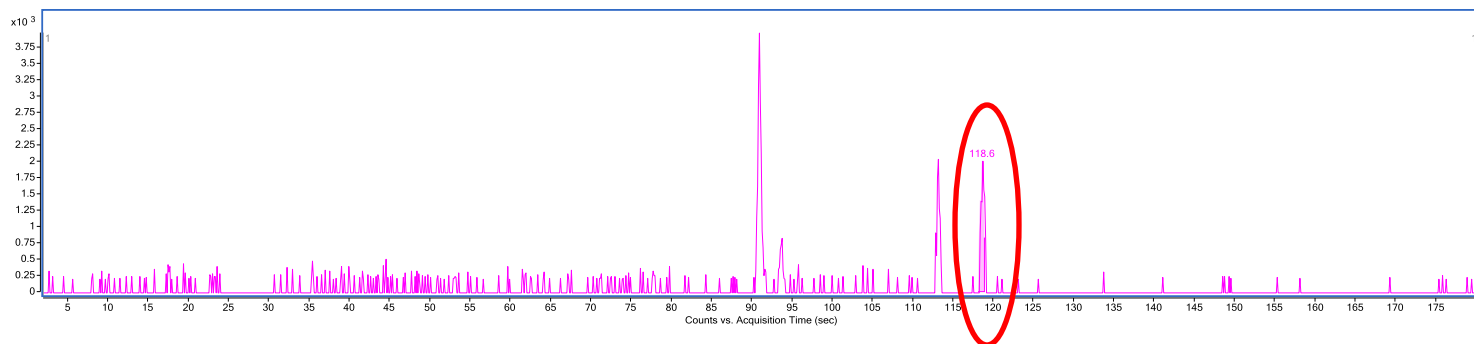
1290 Infinity/6540 Q-TOF

- Tebufenoxide 25pg on-column 3min

MS Formula Results: + Scan [118.6 sec]

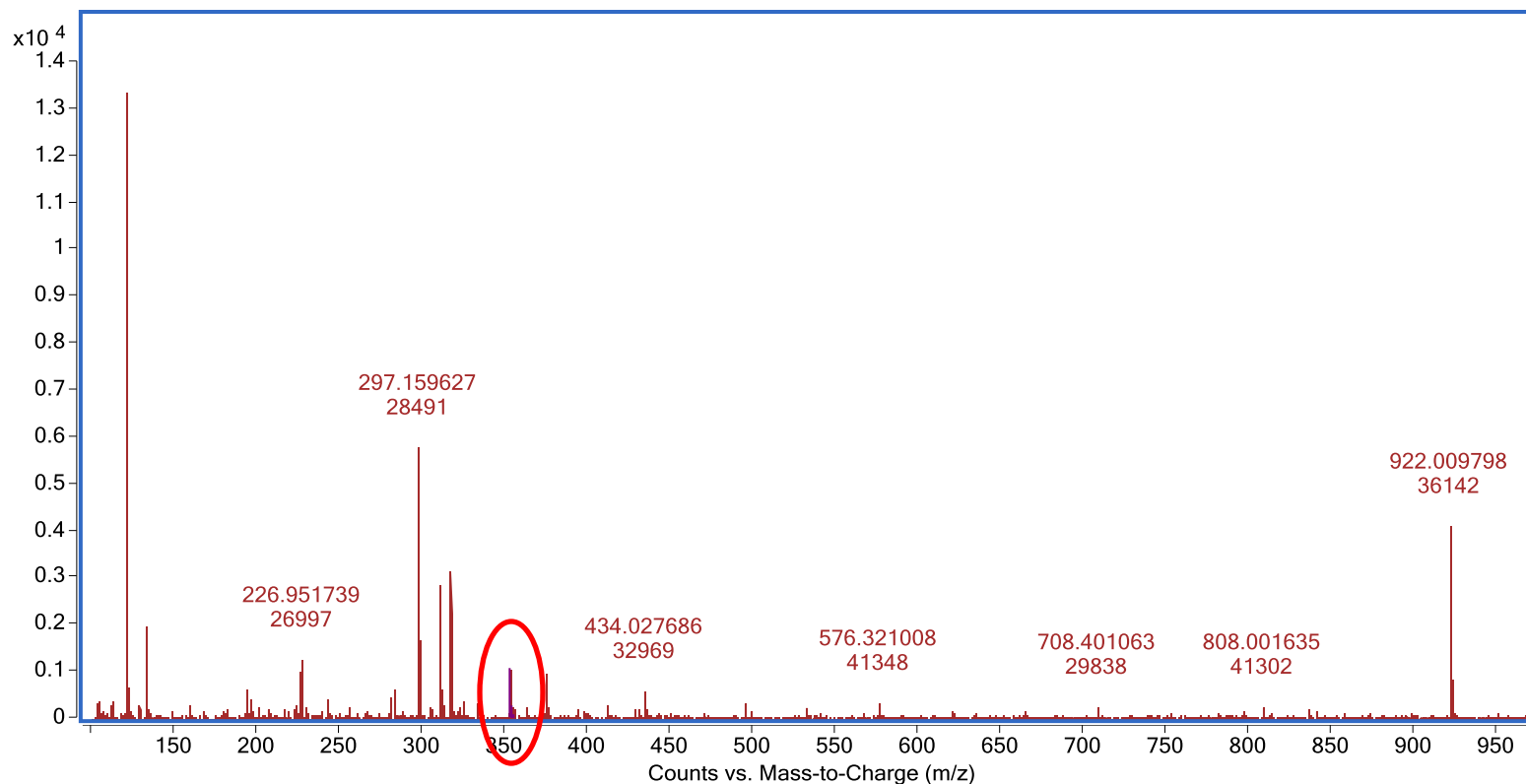
m/z	Ion	Formula	Abundance
353.222598	(M+H)+	C22 H29 N2 O2	1059.5

Best	Formula (M)	Ion Formula	Score	Cross Score	Calc m/z	Diff (ppm)	Mass Match	Abund Match	Spacing Match
☐	C27 H28	C27 H29	64.68		353.226377	10.76	40.34	80.44	94.47
☒	C22 H28 N2 O2	C22 H29 N2 O2	98.56		353.222355	-0.67	99.65	98.34	96.64
☐	C21 H33 Cl O2	C21 H34 Cl O2	68.77		353.224184	4.53	85.11	0.14	84.85
☐	C19 H32 N2 O2 S	C19 H33 N2 O2 S	71.88		353.225725	-8.9	53.76	84.99	52.37
☐	C18 H24 N8	C18 H25 N8	79.33		353.219669	-8.31	58.17	97.62	99.71
☐	C17 H29 Cl N6	C17 H30 Cl N6	64.95		353.221499	-3.11	92.68	0.13	87.27
☐	C16 H33 Cl N2...	C16 H34 Cl N2 O4	53.27		353.220162	-6.89	68.89	0.07	85.88
☐	C15 H28 N8 S	C15 H29 N8 S	91.66		353.22304	1.25	98.79	81.23	89.94
☐	C14 H33 Cl N6 S	C14 H34 Cl N6 S	54.95		353.22487	6.45	72.19	0.02	86.37
☐	C14 H32 N4 O4 S	C14 H33 N4 O4 S	85.5		353.221703	-2.54	95.08	64.55	91.48
☐	C12 H32 N8 S2	C12 H33 N8 S2	50.92		353.226411	10.81	40.04	41.38	84.12
☐	C11 H28 N8 O5	C11 H29 N8 O5	67.49		353.225543	8.35	57.91	56.88	99.39



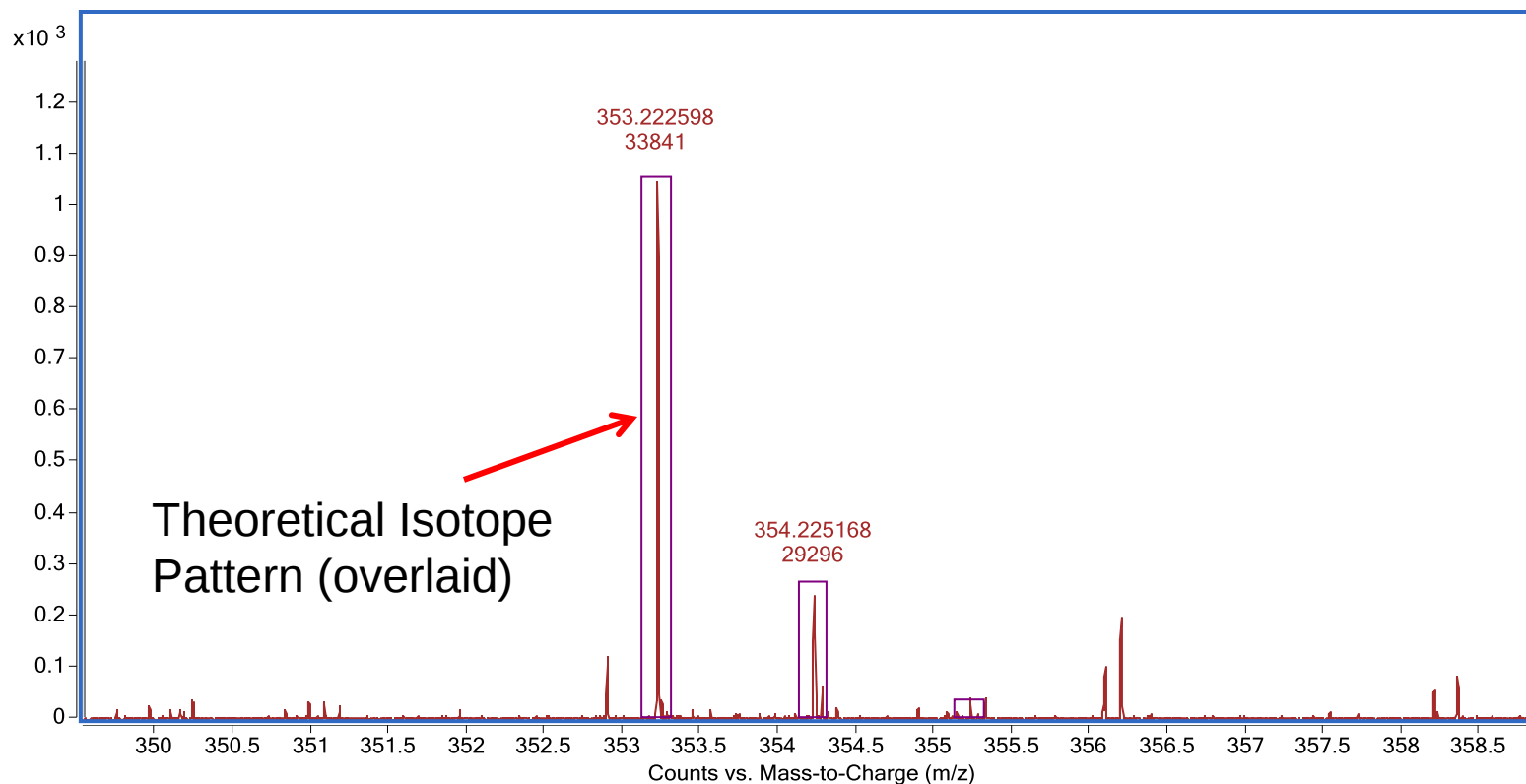
1290 Infinity/6540 Q-TOF

- Tebufenoxide 25pg on-column 3min



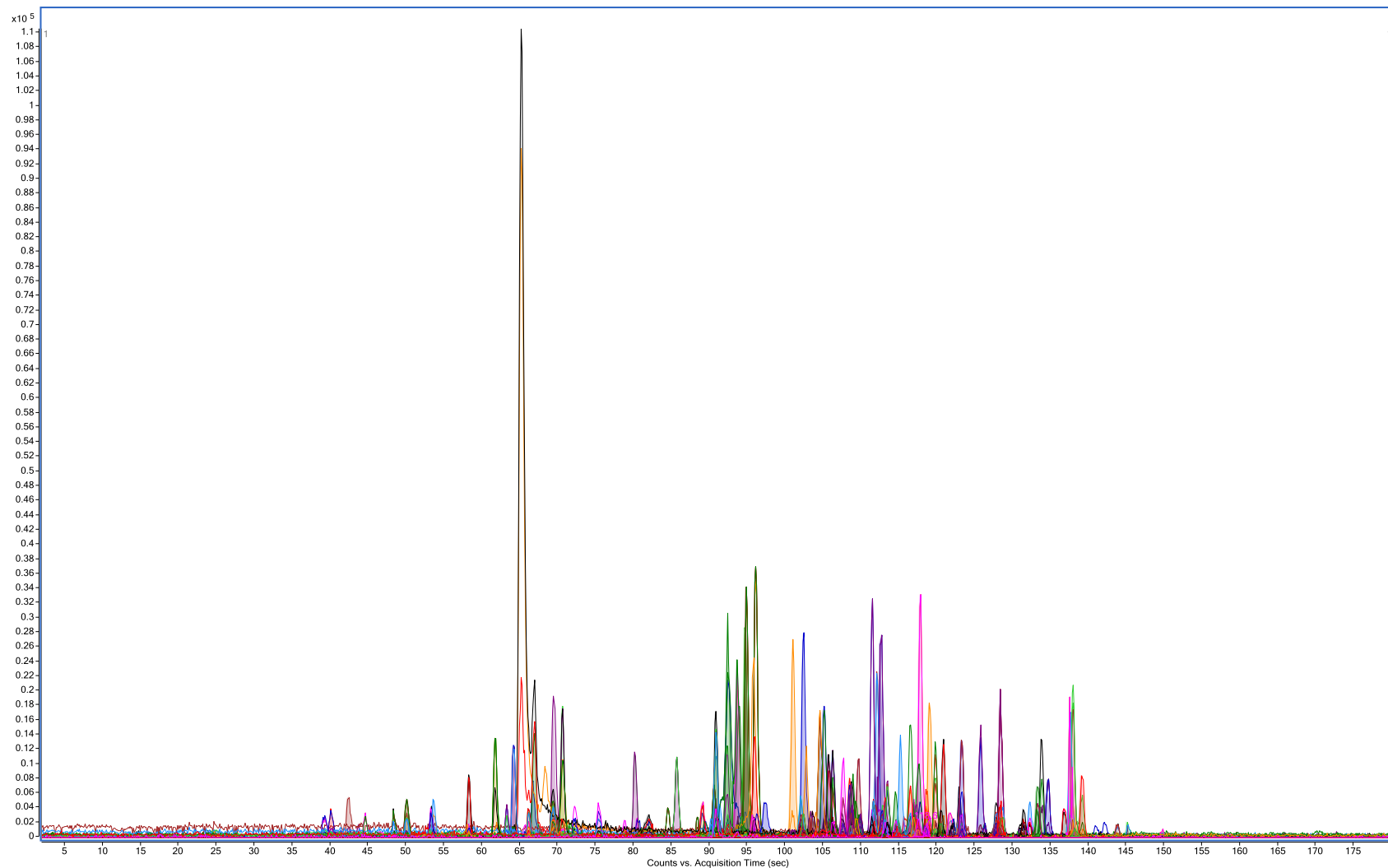
1290 Infinity/6540 Q-TOF

- Tebufenoxide 25pg on-column 3min



25pg on-column 1290 Infinity/6540 Q-TOF

3min 92% identified pesticides



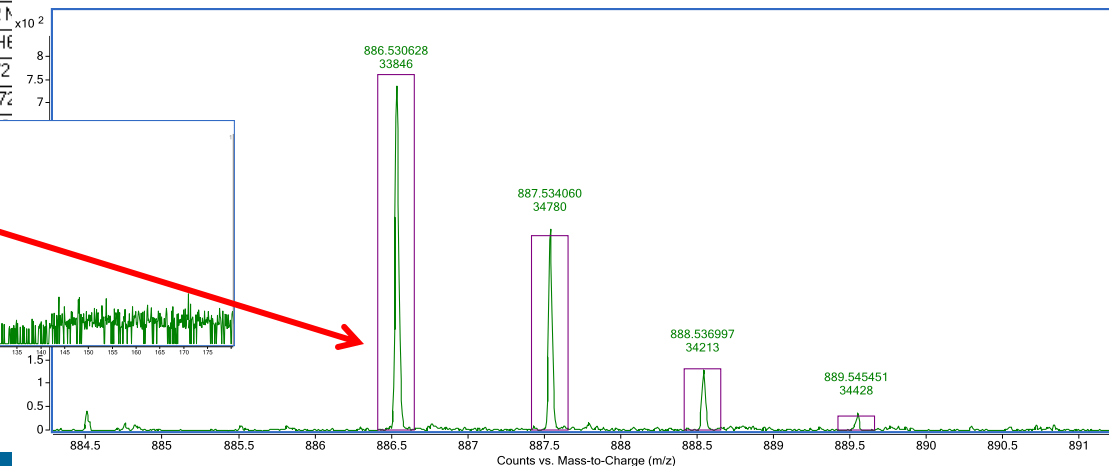
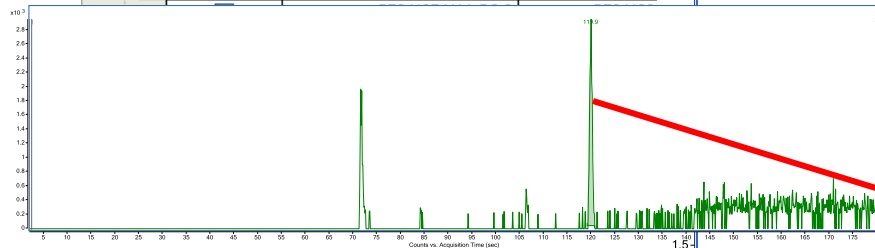
1290 Infinity/6540 Q-TOF

- Emamectin B1a, 50pg on-column 3min

m/z	Ion	Formula	Abundance
886.530628	(M+H)+	C49 H76 N 013	737.4

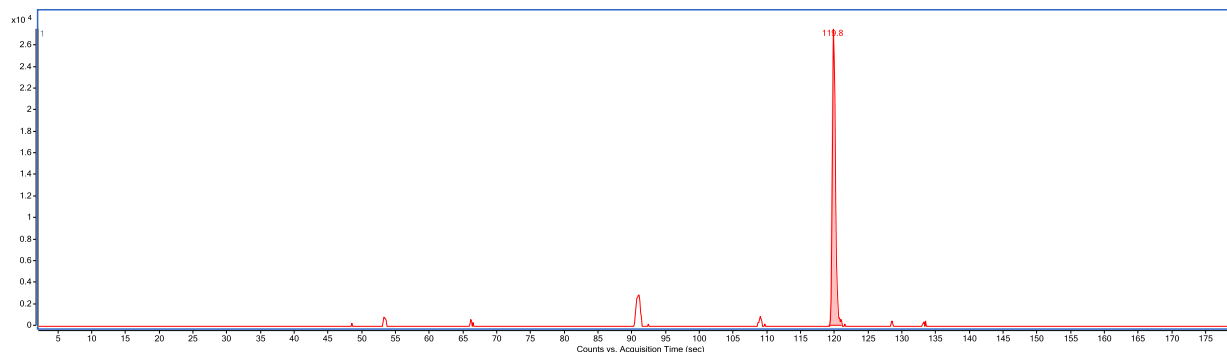
Best	Formula (M)	Ion Formula	Score	Calc m/z	Diff (ppm)	Mass Match	Abund Match	Spacing Match
☐	C47 H63 N15 O3	C47 H64 N15 O3	98.97	886.531107	0.41	99.72	97.9	98.77
☒	C49 H75 N 013	C49 H76 N 013	98.59	886.531118	0.54	99.53	95.87	99.99
☐	C46 H67 N11 O7	C46 H68 N11 O7	97.89	886.52977	-1.07	98.18	96.24	99.28
☐	C50 H71 N5 O9	C50 H72 N5 O9	96.52	886.532455	2.02	93.63	99.54	99.88
☐	C54 H71 N5 O4 S	C54 H72 N5 O4 S	94.76	886.529953	-0.83	98.9	86.14	96.8
☐	C55 H67 N9 S	C55 H68 N9 S	94.63	886.53129	0.66	99.3	85.44	96.31
☐	C40 H63 N21 O 5	C40 H64 N21 O 5	93.07	886.531793	1.1	98.06	88.22	88.9
☐	C45 H71 N7 O11	C45 H72 N7 O11	92.9	886.528433	-2.55	90.05	92.01	99.66
☐	C43 H59 N21 O	C43 H60 N21 O	92.56	886.528422	-2.67	89.08	94.48	97.2
☐	C39 H67 N17 O5 S	C39 H68 N17 O5 S	91.85	886.530456	-0.38	99.77	80.29	89.91
☐	C53 H75 N 08 S	C53 H76 N 08 S	91.13	886.528616	-2.31	91.75	85.03	97.19
☐	C51 H67 N9 O5	C51 H68 N9 O5	91	886.533793	3.5	82.05	98.73	99.64
☐	C42 H75 N7 O11 S	C42 H76 N7 O11 S	90.28	886.531803	1.23	97.57	75.25	93.73
☐	C58 H63 N9	C58 H64 N9	89.72	886.52792	-3.13	85.37	88.62	99.75
☐	C41 H79 N3 O15 S	C41 H80 N3 O15 S	88.73	886.530466	-0.24	99.91	65.48	94.28
☐	C43 H71 N11 O7 S	C43 H72 N11 O7 S	88.31	886.533141	2.71	88.79	83.55	93.05
☐	C47 H71 N11 O2 S2	C47 H72 N11 O2 S2						
☐	C32 H59 N27 O4	C32 H60 N27 O4						
☐	C38 H71 N13 O9 S	C38 H72 N13 O9 S						
☐	C34 H71 N13 O14	C34 H72 N13 O14						

886.530628
33846

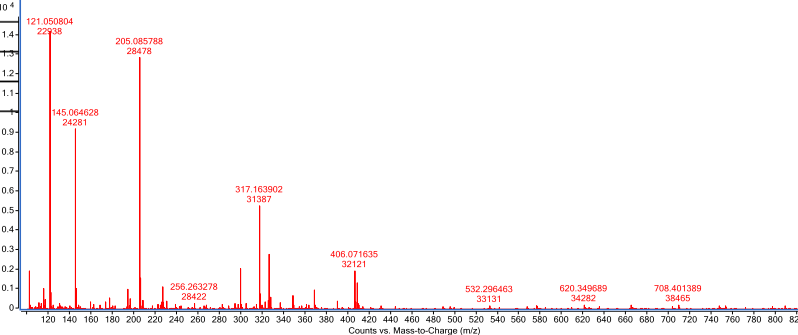
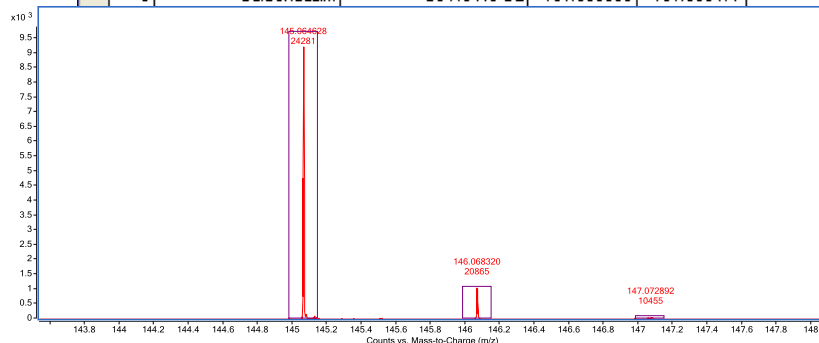


1290 Infinity/6540 Q-TOF

- 1-naphthol, 50pg on-column 3min

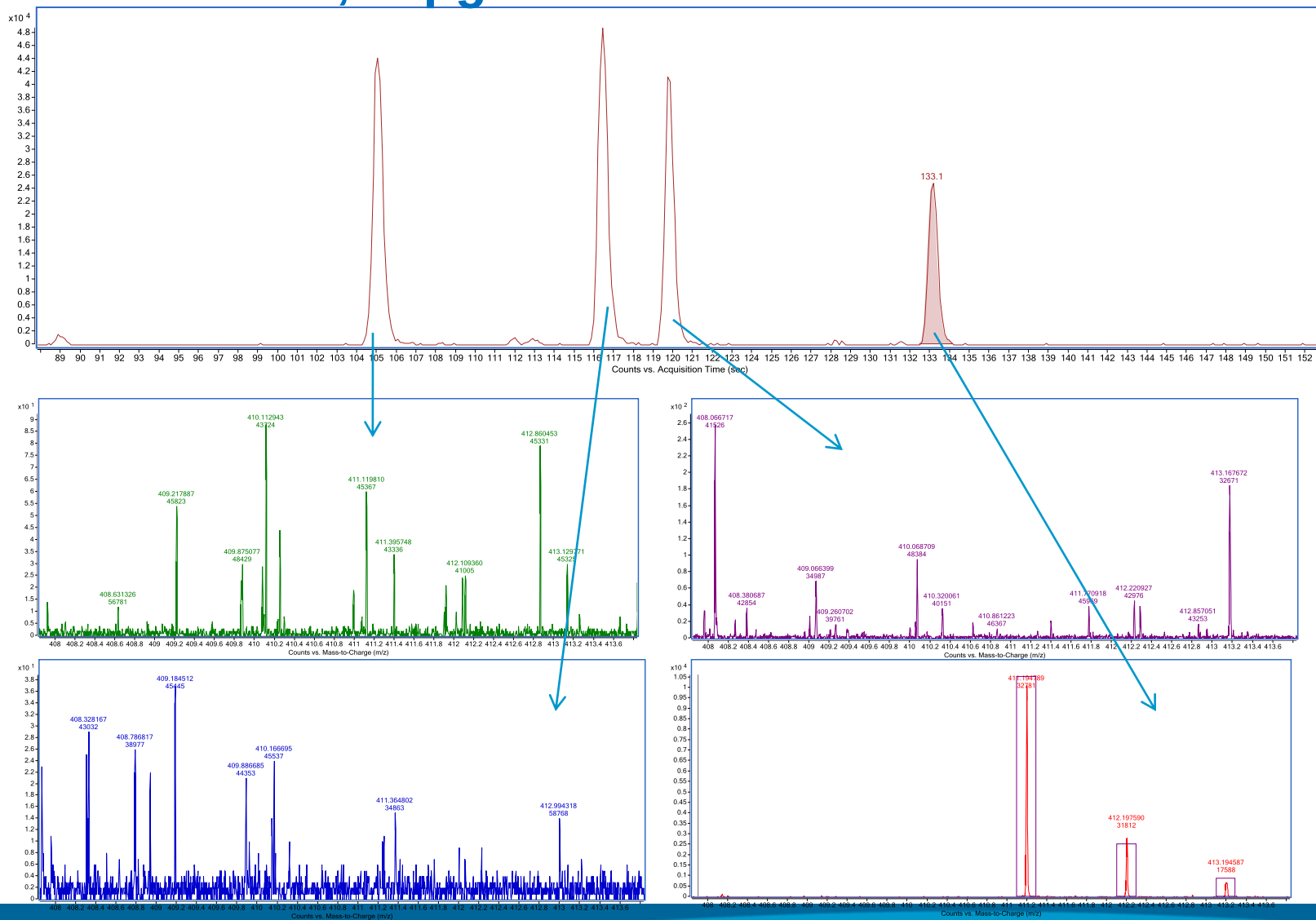


Compound List												
Cpd	Name	Formula (Tgt)	Mass	Mass (Tgt)	Diff (Tgt, ppm)	RT	Score (Tgt)	Height	Area	Abund	Algorithm	
1	Methamidophos	C2 H8 N O2 P S	141.001262	141.001336	-0.52	24.6	47.59	1590	639	823	Find By Formula	
2	Naphthol, 1-	C10 H8 O	144.057502	144.057515	-0.09	119.8	99.16	27504	14953	15022	Find By Formula	
3	Methomyl	C5 H10 N2 O2 S	162.04672	162.046298	2.6	92.9	46.69	559	506	296	Find By Formula	
4	Fenuron	C9 H12 N2 O	164.094596	164.094963	-2.24	66.8	98.56	16508	11151	7436	Find By Formula	
5	Fuberidazole	C11 H8 N2 O	184.063654	184.063663	-0.05	69.5	99.9	53636	31874	31498	Find By Formula	
6	Propamocarb	C9 H20 N2 O2	188.15212	188.152478	-1.9	42.3	99.27	14920	10513	7714	Find By Formula	
7	Tricyclazole	C9 H7 N3 S	189.035792	189.036068	-1.46	70.6	104.4					
8	Butocarboxim	C7 H14 N2 O2 S	190.077627	190.077598	0.15	78.4	14.4					
9	Carbendazim	C9 H9 N3 O2	191.069335	191.069477	-0.74	61.7	12.7					



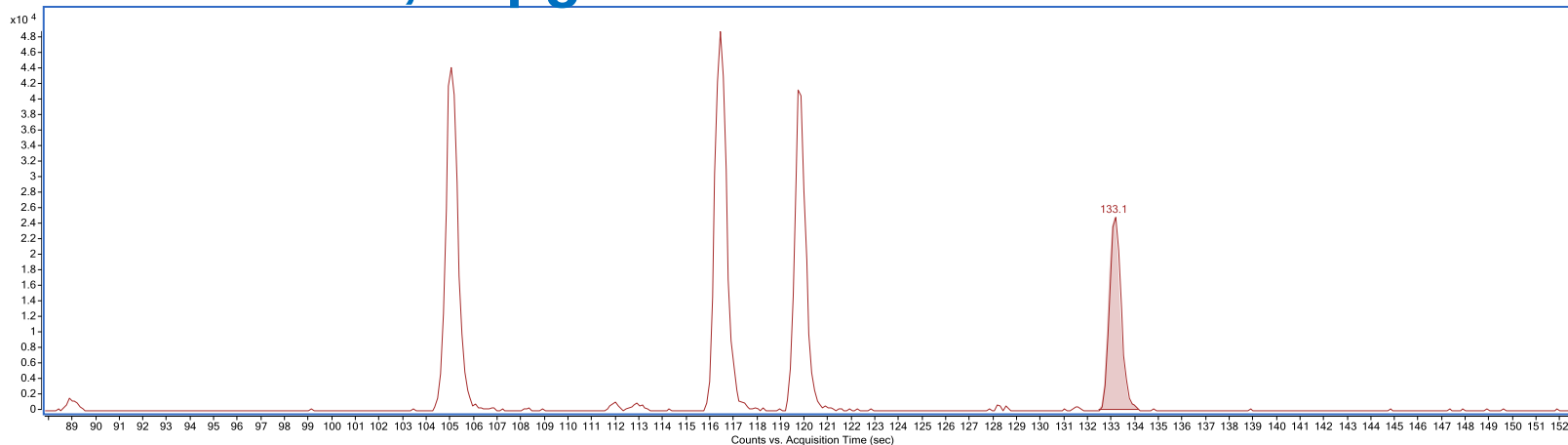
1290 Infinity/6540 Q-TOF

- Benfuracarb, 50pg on-column 3min



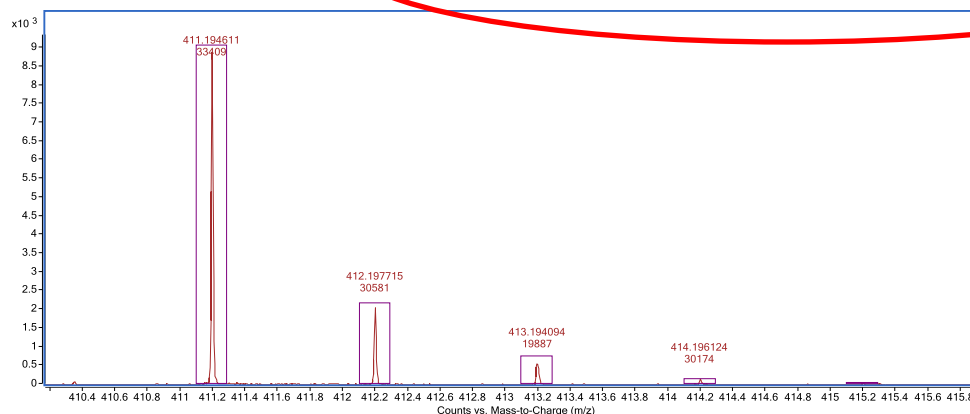
1290 Infinity/6540 Q-TOF

- Benfuracarb, 50pg on-column 3min



MS Formula Results: + Scan (132.9-133.4 sec)

m/z	Ion	Formula	Abundance						
411.194611	(M+H)+	C20 H31 N2 O5 S	9225						
Best	Formula (M)	Ion Formula	Score	Calc m/z	Diff (ppm)	Mass Match	Abund Match	Spacing Match	
☒	C20 H30 N2 O5 S	C20 H31 N2 O5 S	98.62	411.194819	0.51	99.77	95.92	99.56	



Non-Targeted Screen Summary (QTOF/TOF)

- Agilent TOF/Q-TOF for 'unknown' screening (low ppm mass accuracy.)
- Agilent MassHunter PCD utilizes the accurate mass data to provide reliable 'positives' for qualitative screens.
 - Extra Confirmation with Isotope abundance (pattern), spacing and mass matching.
 - 1600 Pesticide analytes
 - Fully Automated integration with MassHunter Qualitative Analysis SW.
- Reliable Data mining Algorithms (MFE or FBF.)
- Low pico-gram on-column Sensitivity.
- Fully Automated.
- Faster data acquisition with 1290 Infinity UHPLC.

Questions?

Thank You for Listening.



Agilent Technologies