

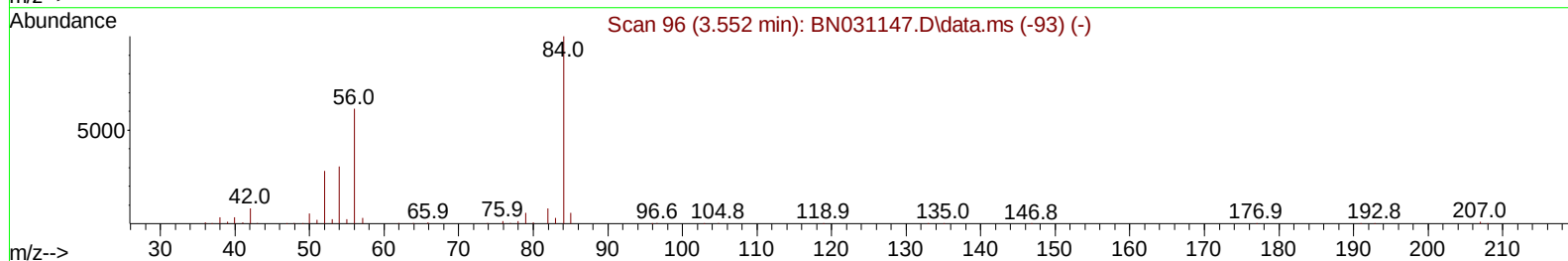
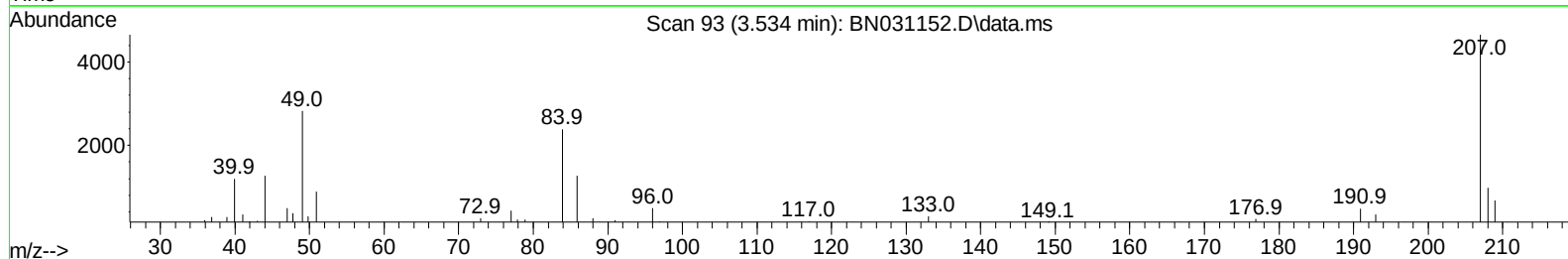
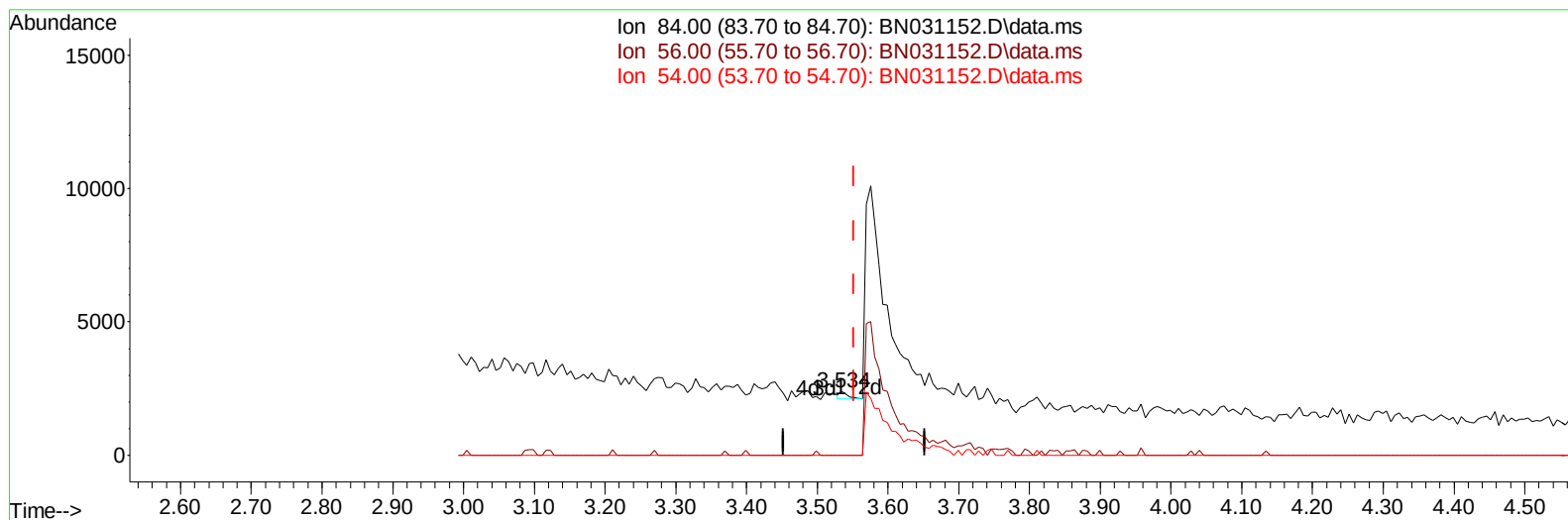
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042224\
 Data File : BN031152.D
 Acq On : 23 Apr 2024 12:58
 Operator : MA/JU
 Sample : P2159-12DL 5X
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORD3DL

Manual IntegrationsAPPROVED

Quant Time: Apr 23 22:40:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 19 22:56:25 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/24/2024
 Supervised By :mohammad ahmed 04/24/2024



TIC: BN031152.D\data.ms

(4) Pyridine-d5 (S)

3.534min (-0.018) 0.03 ng/u1

response 225

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	0.00#
54.00	30.10	0.00#
0.00	0.00	0.00

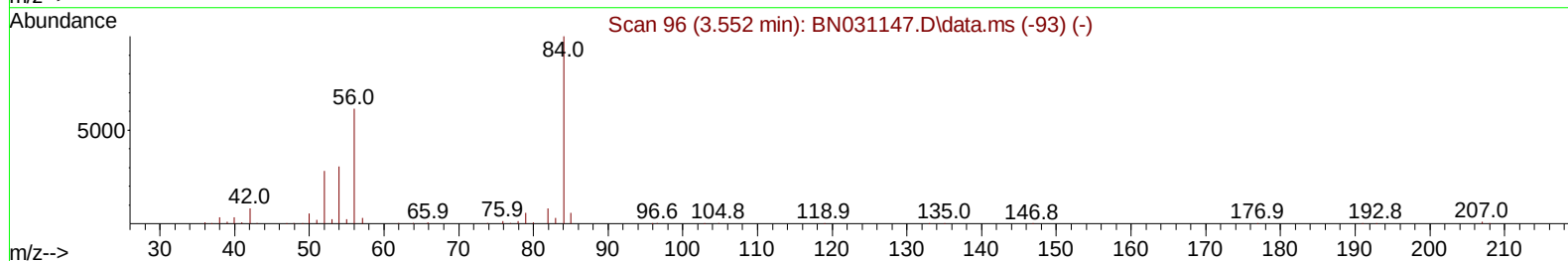
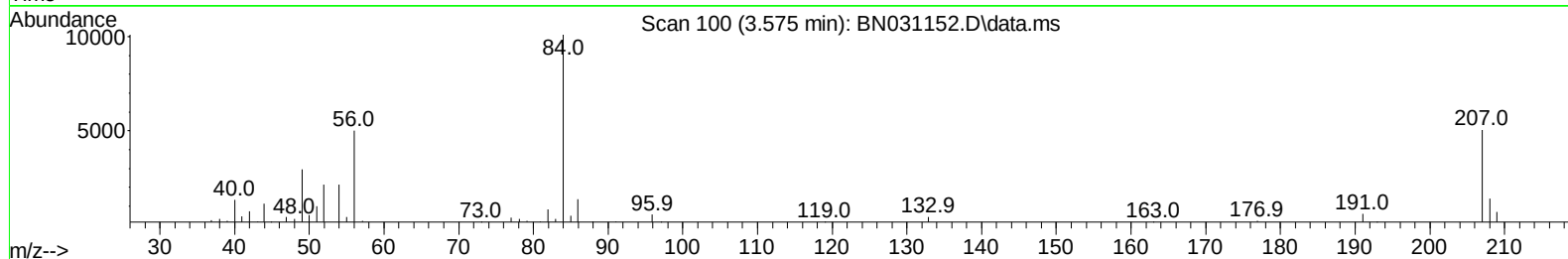
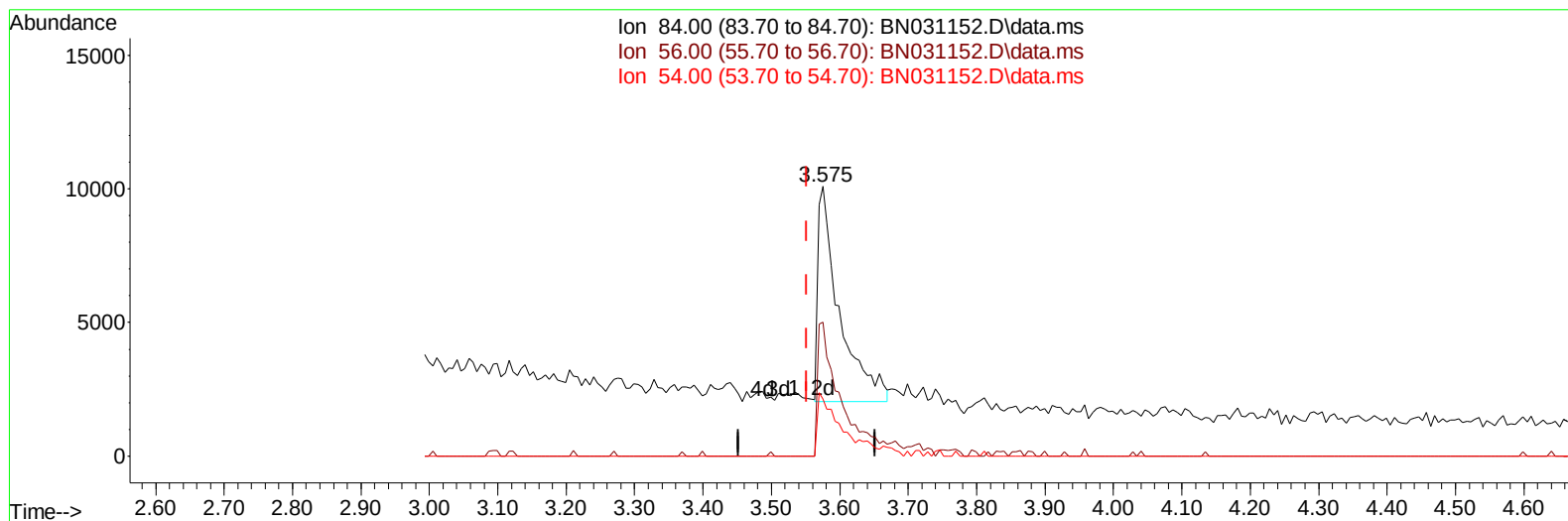
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(4) Pyridine-d5 (S)

3.575min (+ 0.023) 2.20 ng/ul m

response 17615

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	49.56
54.00	30.10	21.11#
0.00	0.00	0.00

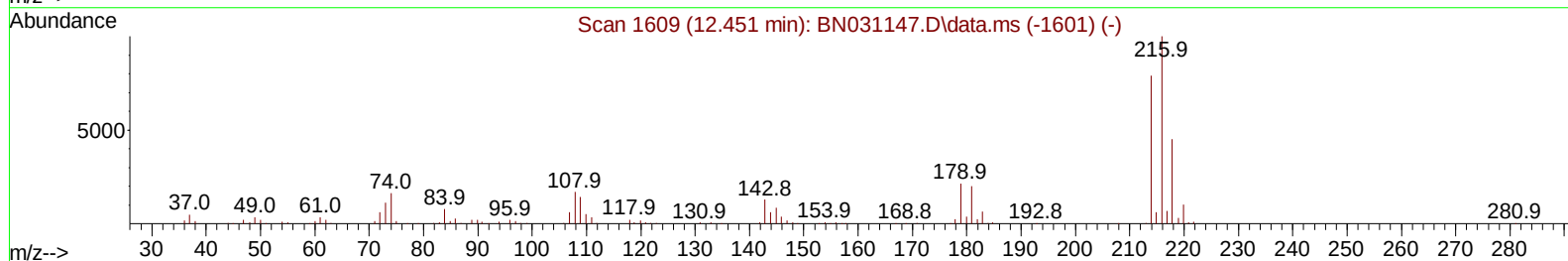
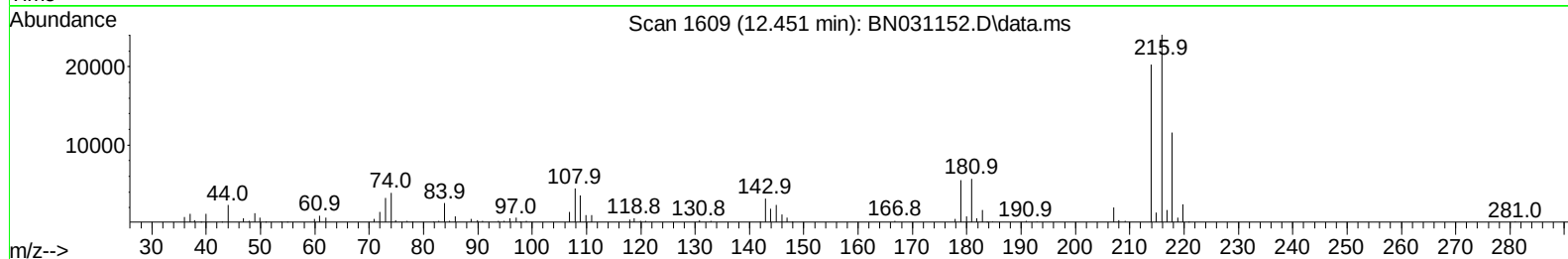
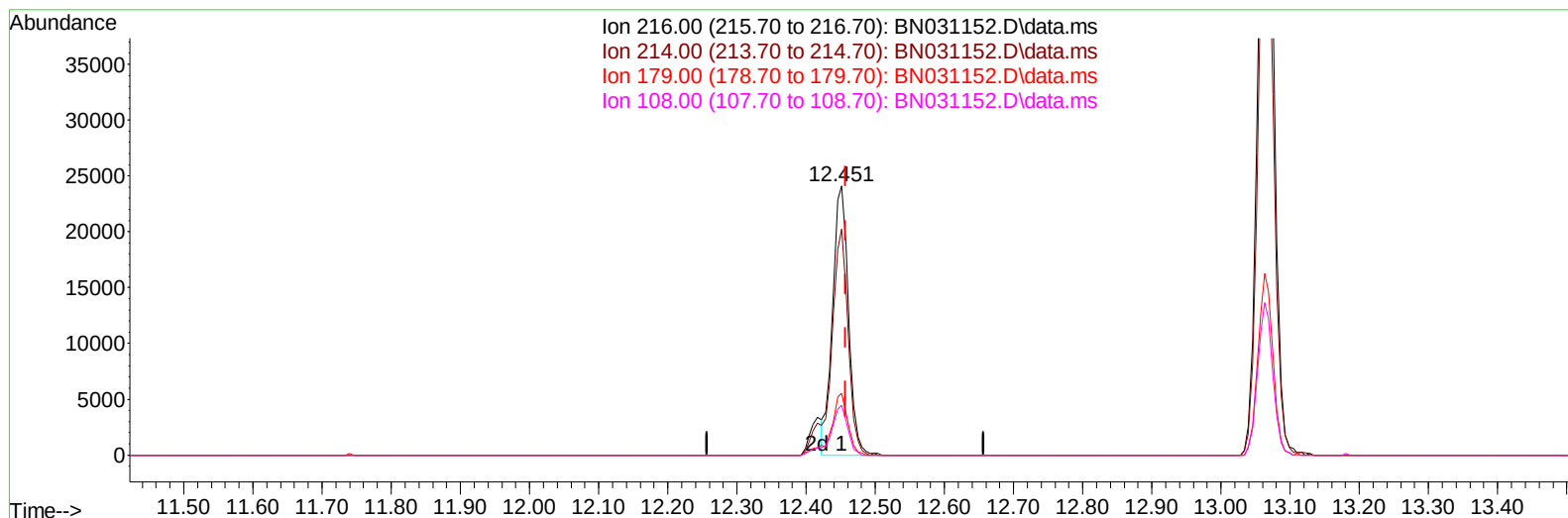
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TIC: BN031152.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.451min (-0.006) 3.48 ng/uI

response 38434

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	78.90	84.05
179.00	21.30	23.04
108.00	17.80	18.63

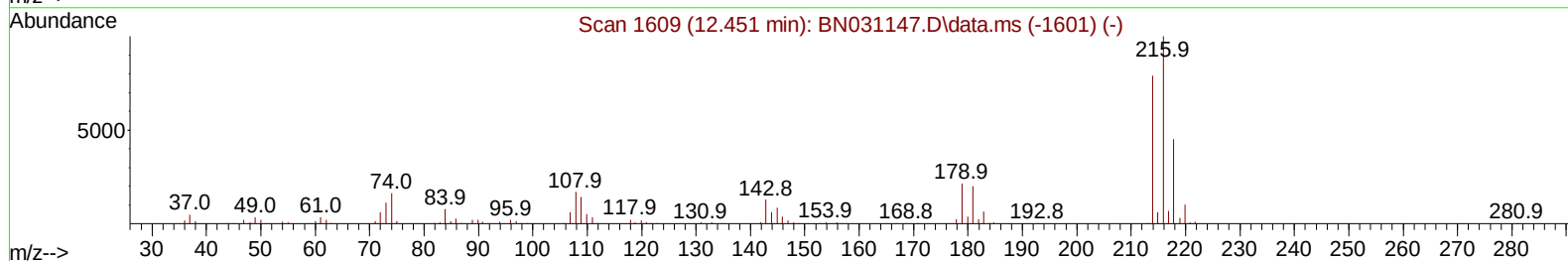
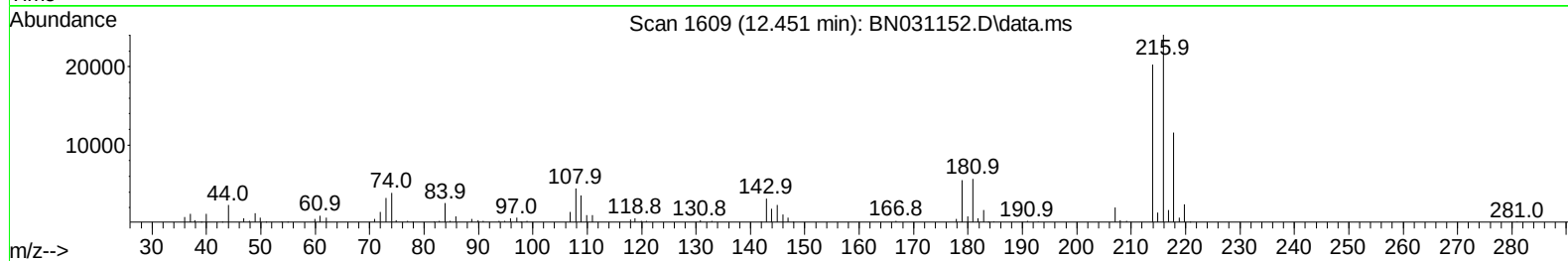
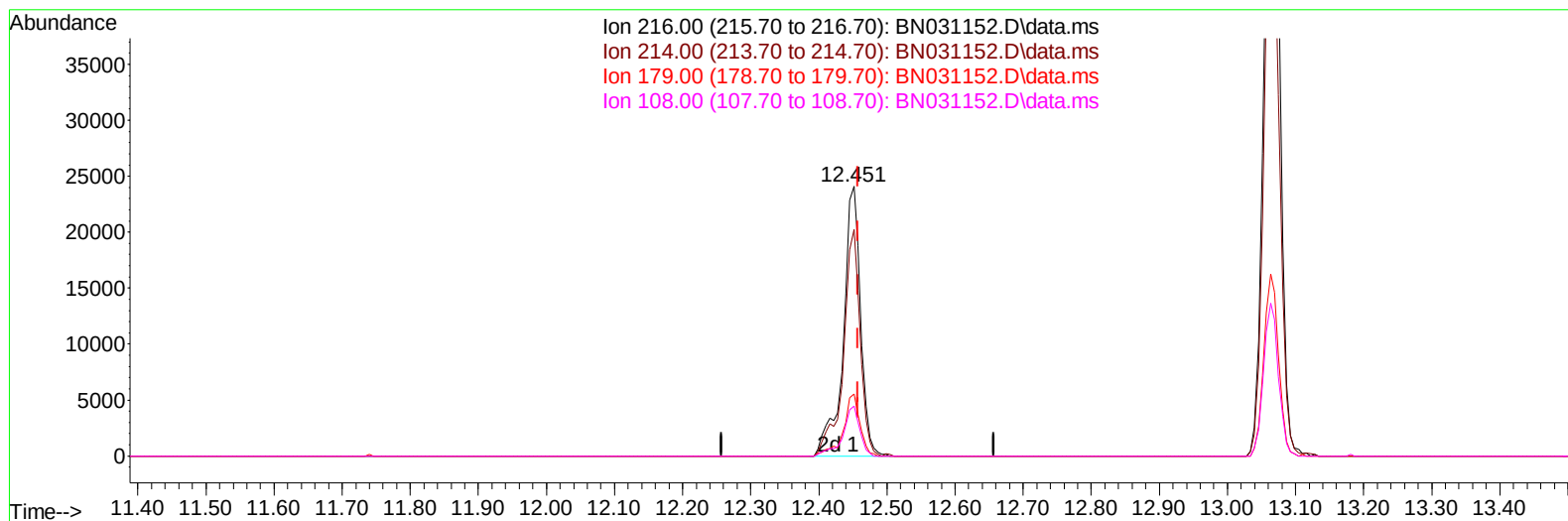
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ALS Vial : 41 Sample Multiplier: 1

Instrument :
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TIC: BN031152.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.451min (-0.006) 3.86 ng/ul m

response 42559

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	78.90	84.05
179.00	21.30	23.04
108.00	17.80	18.63

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/24/2024
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Quant Time: Apr 23 23:05:16 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	129651	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	602924	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164	421611	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	936295	20.000	ng/ul	0.00
79) Chrysene-d12	21.263	240	988200	20.000	ng/ul	0.00
88) Perylene-d12	24.157	264	1100496	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	2255	0.813	ng/uL	0.00
4) Pyridine-d5	3.575	84	17615m	2.204	ng/ul	0.02
7) Phenol-d5	6.811	99	11654	1.066	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	43638	6.858	ng/ul	0.00
11) 2-Chlorophenol-d4	7.163	132	40610	4.769	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	25287	2.749	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	27159	6.269	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	24976	5.547	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	48699	5.484	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	65629	4.729	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	207910	7.087	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	224756	6.760	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.275	176	185746	7.273	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	16385	3.469	ng/ul	0.00
73) Anthracene-d10	17.128	188	303893	7.488	ng/ul	0.00
81) Pyrene-d10	19.427	212	389423	8.019	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	376730	7.219	ng/ul	-0.02
Target Compounds						
					Qvalue	
22) Nitrobenzene	8.828	77	226074	22.112	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.451	216	42559m	3.858	ng/ul	
75) 1,2,3,4-Tetrachloroben...	13.063	216	105591	9.615	ng/uL	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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