

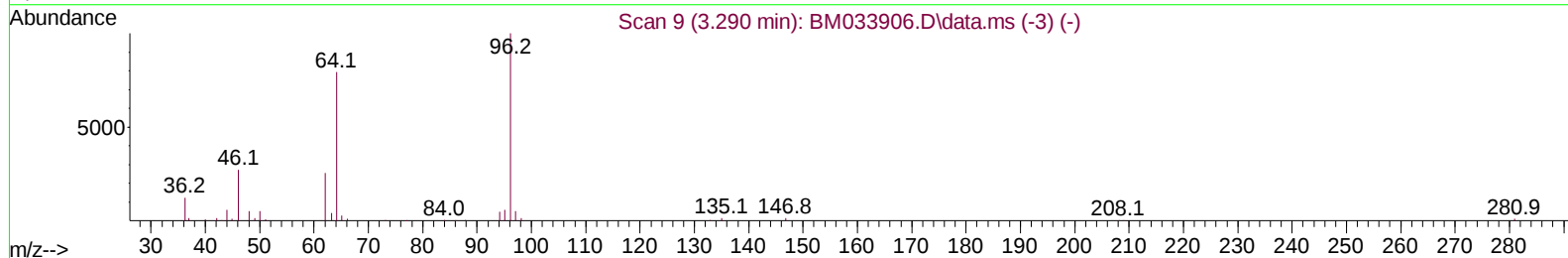
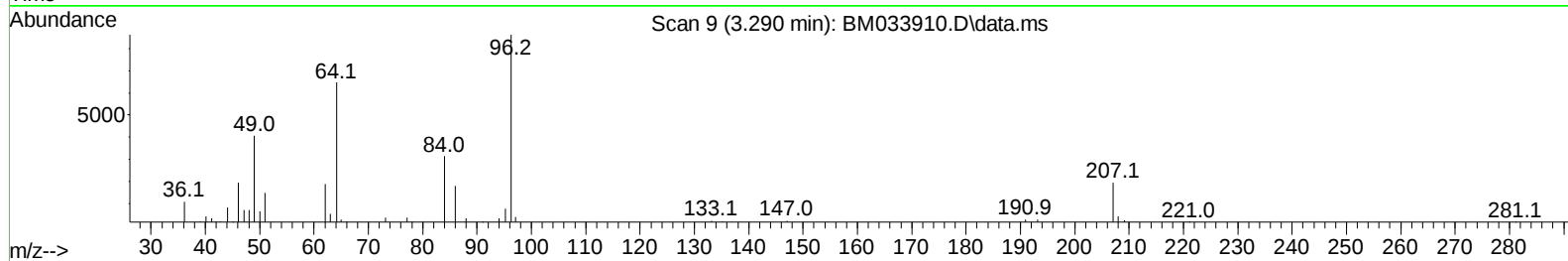
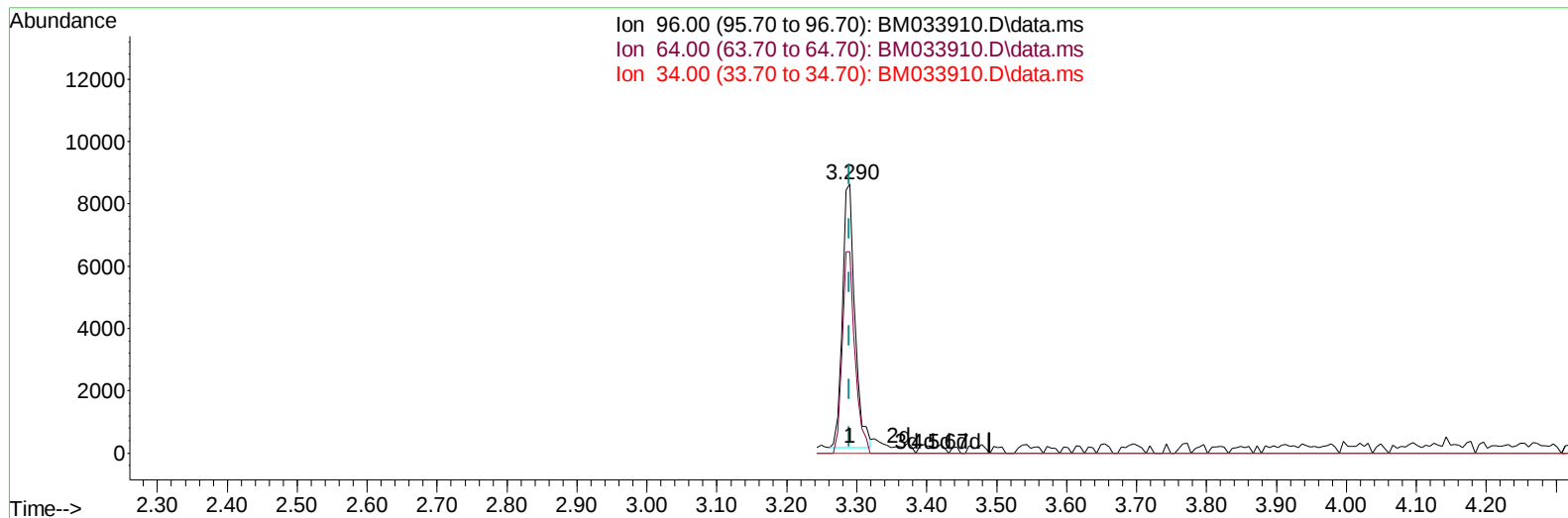
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033910.D
 Acq On : 28 Jan 2022 18:55
 Operator : CG/JU
 Sample : N1302-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F53

Manual IntegrationsAPPROVED

Quant Time: Jan 28 23:28:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 28 23:26:29 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2022
 Supervised By :mohammad ahmed 02/01/2022



TIC: BM033910.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.290min (+ 0.000) 4.06 ng/uL

response 10647

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	74.98
34.00	0.00	0.00
0.00	0.00	0.00

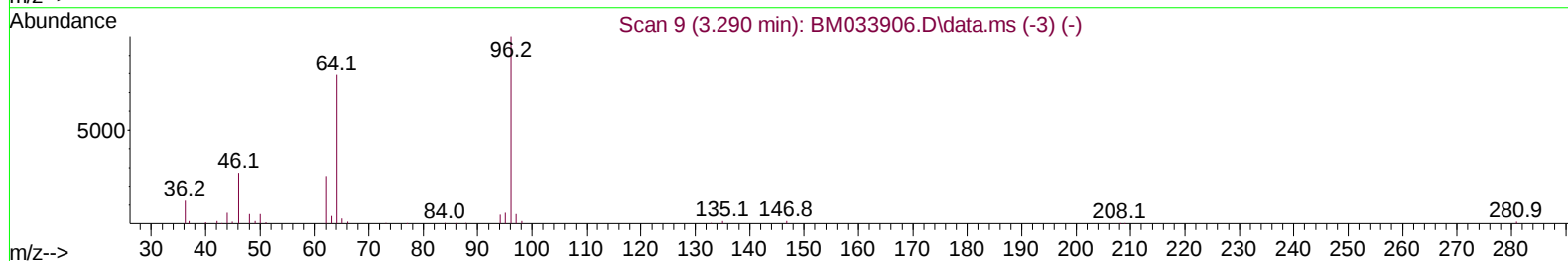
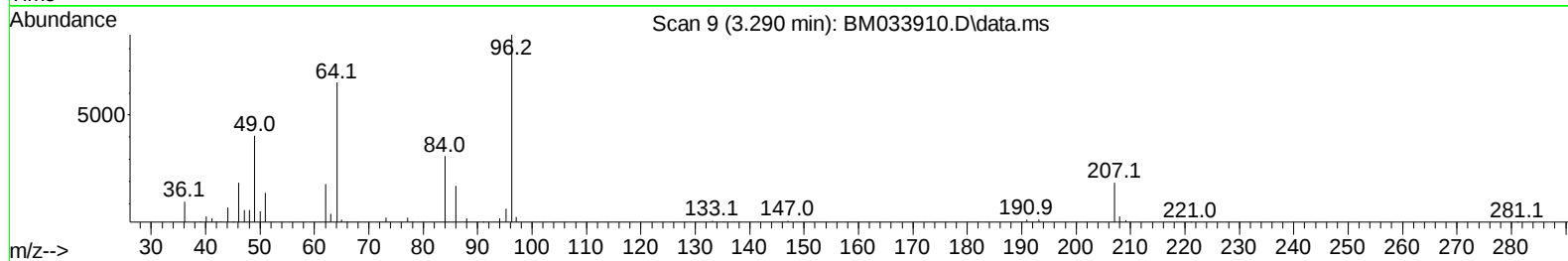
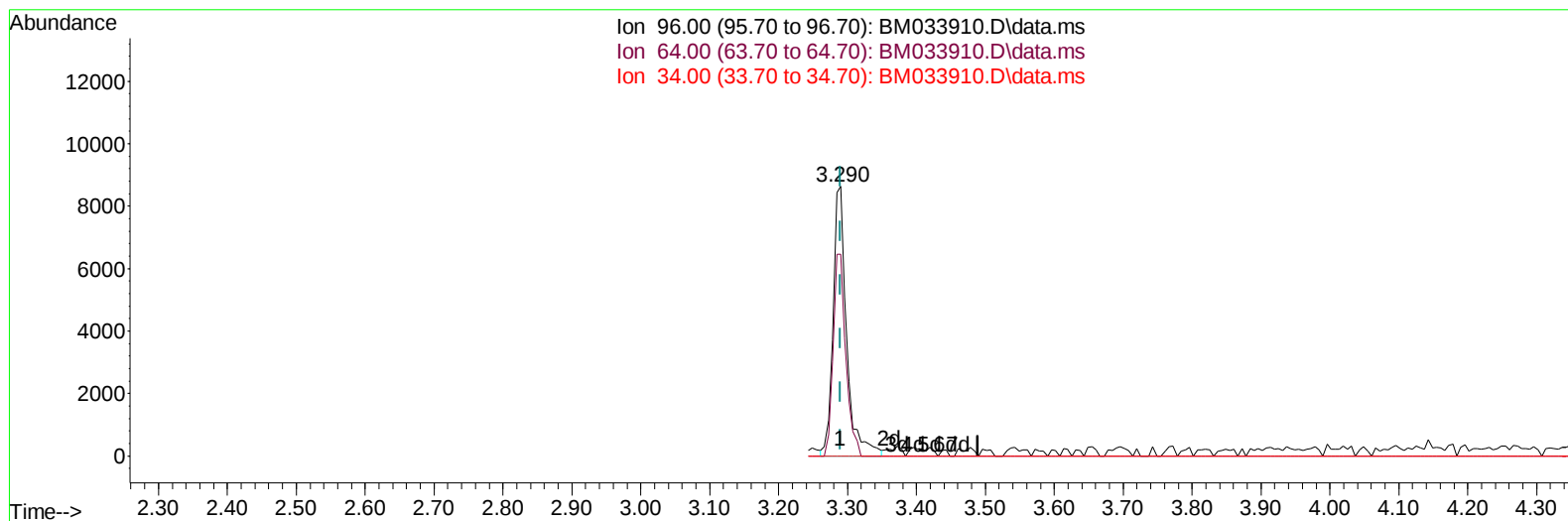
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TIC: BM033910.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.290min (+ 0.000) 4.49 ng/uL m

response 11791

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	74.98
34.00	0.00	0.00
0.00	0.00	0.00

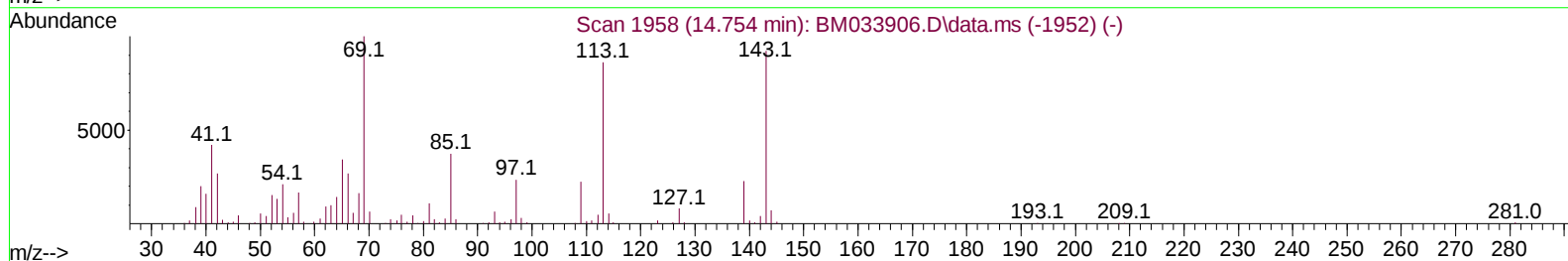
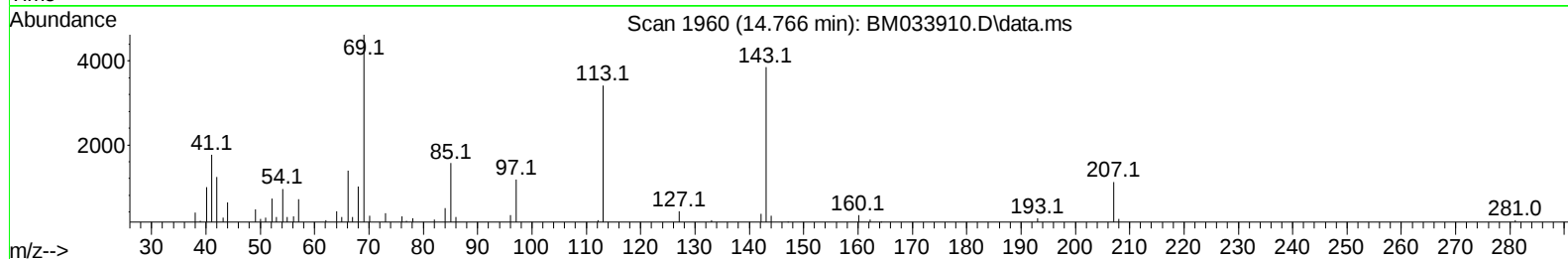
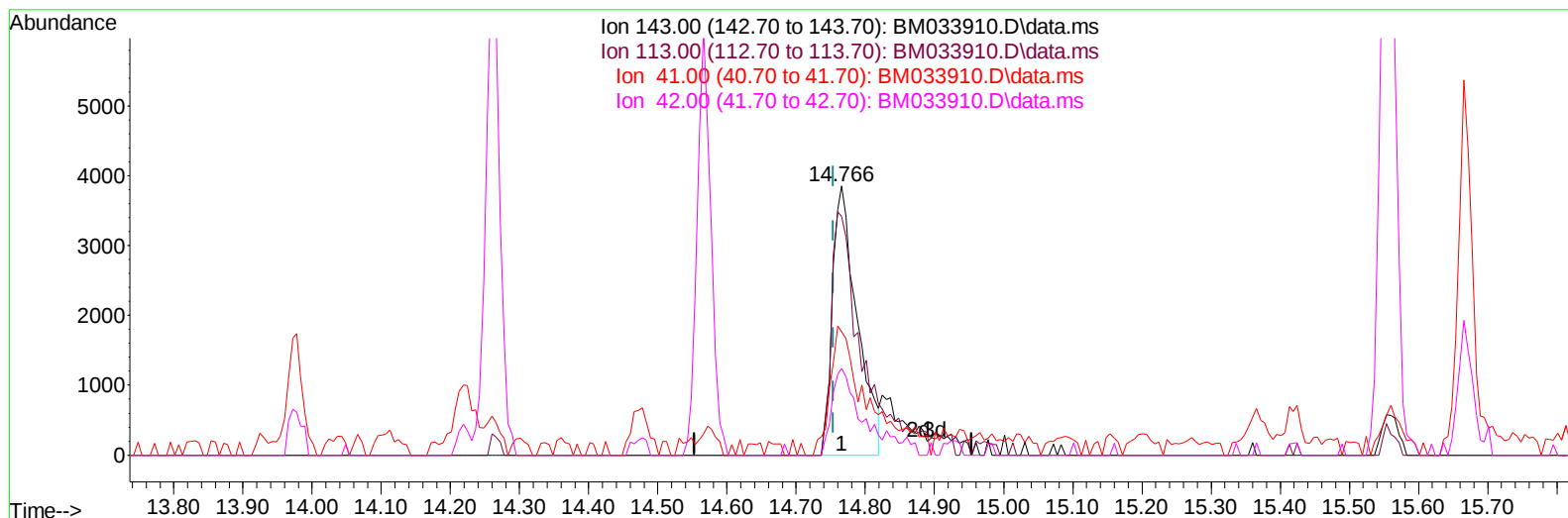
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TIC: BM033910.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 2.58 ng/u1

response 9529

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	88.81
41.00	57.20	45.88
42.00	39.50	32.15

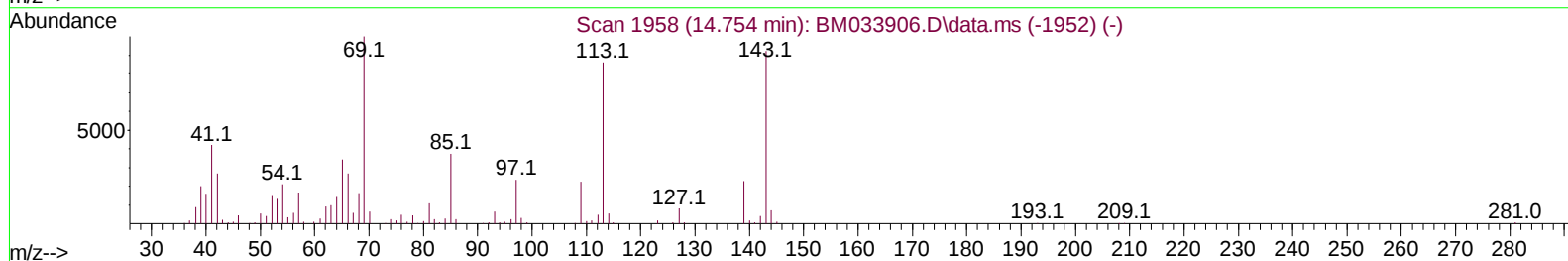
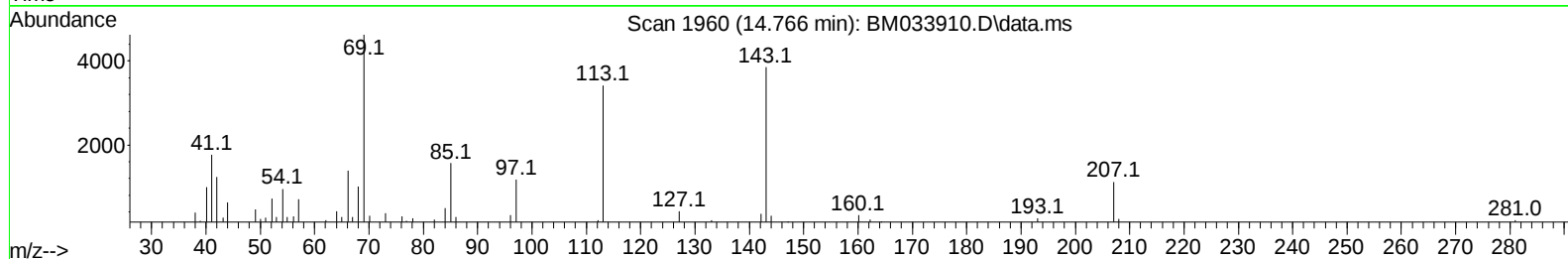
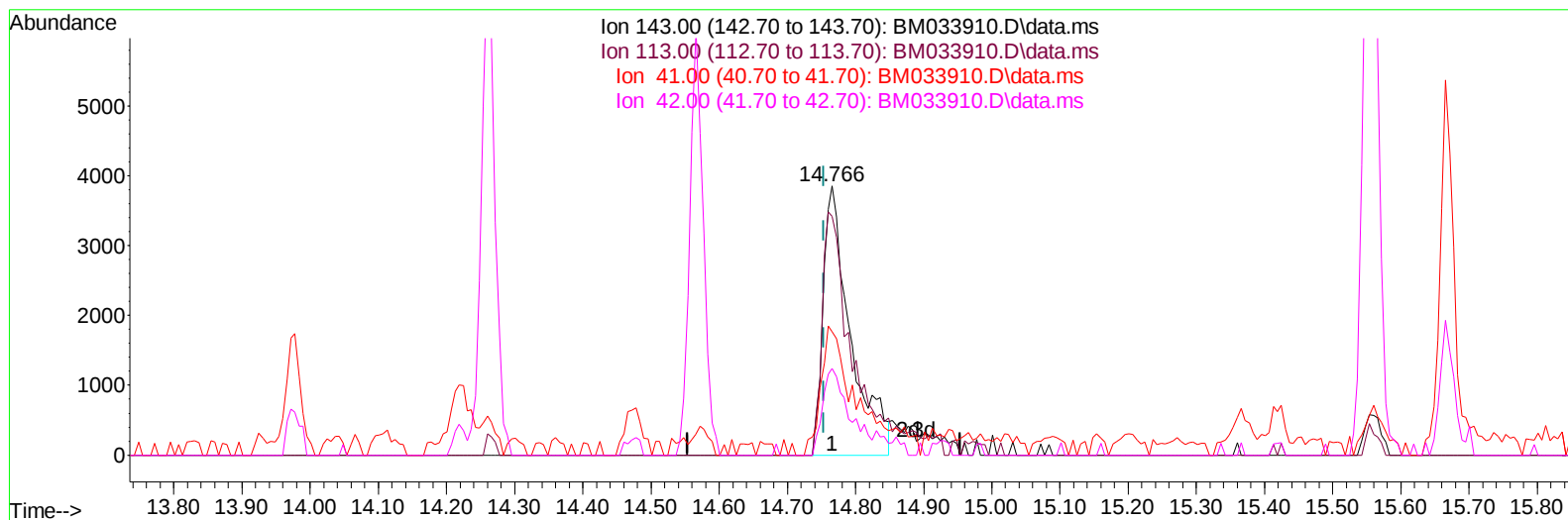
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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TIC: BM033910.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 2.91 ng/ul m

response 10750

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	88.81
41.00	57.20	45.88
42.00	39.50	32.15

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	98572	20.000	ng/ul	0.00
20) Naphthalene-d8	10.743	136	402704	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	244568	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.307	188	489504	20.000	ng/ul	0.00
79) Chrysene-d12	21.471	240	464227	20.000	ng/ul	0.00
88) Perylene-d12	23.865	264	477110	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	11791m	4.494	ng/uL	0.00
4) Pyridine-d5	3.719	84	53888	7.396	ng/ul	0.00
7) Phenol-d5	7.084	99	52796	5.963	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.255	67	167651	31.096	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	167739	24.781	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	99513	13.578	ng/ul	0.00
21) Nitrobenzene-d5	9.090	128	102659	32.280	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	106300	31.538	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	171932	26.608	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	233129	24.272	ng/ul	0.00
46) Dimethylphthalate-d6	13.978	166	656209	34.055	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	755541	33.236	ng/ul	0.00
54) 4-Nitrophenol-d4	14.766	143	10750m	2.907	ng/ul	0.01
60) Fluorene-d10	15.554	176	544643	34.070	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	78231	26.867	ng/ul	0.00
73) Anthracene-d10	17.401	188	899132	37.854	ng/ul	0.00
81) Pyrene-d10	19.689	212	1016455	40.545	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.712	264	968448	39.012	ng/ul	0.00

Target Compounds Qvalue

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

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