

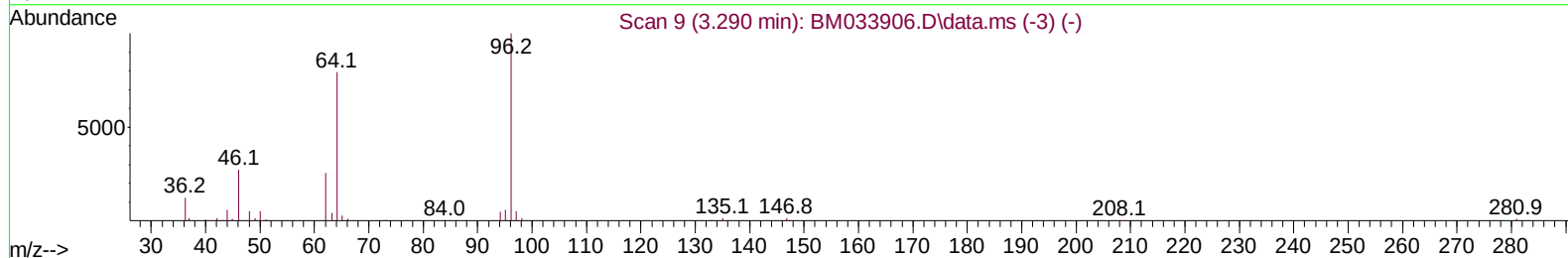
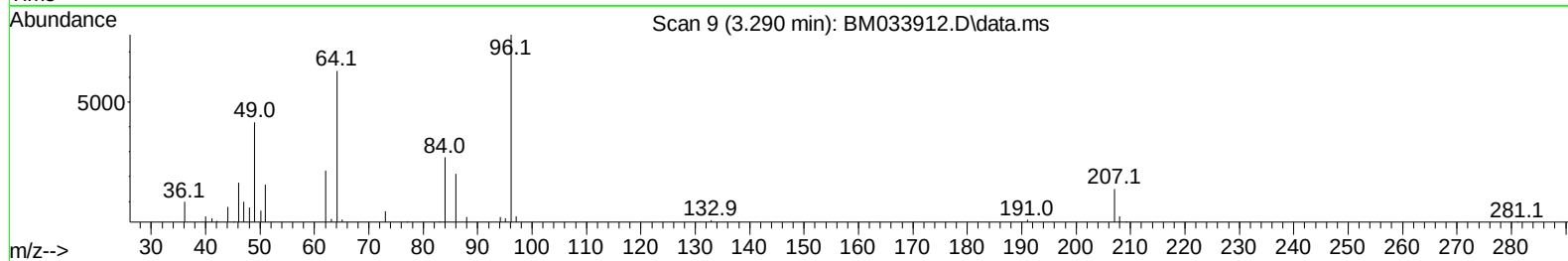
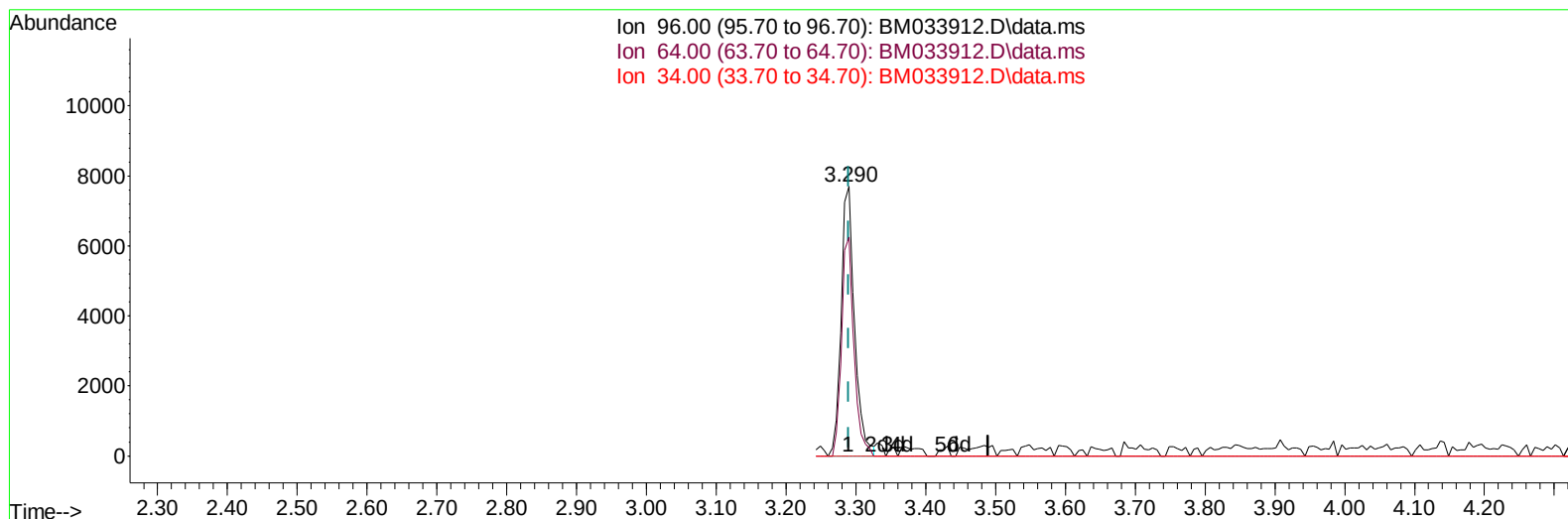
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033912.D
 Acq On : 28 Jan 2022 20:06
 Operator : CG/JU
 Sample : N1302-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F28

Manual IntegrationsAPPROVED

Quant Time: Jan 28 23:29:02 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: BM033912.D\data.ms

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

3.290min (+ 0.000) 4.57 ng/uL

response 10199

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	81.17
34.00	0.00	0.00
0.00	0.00	0.00

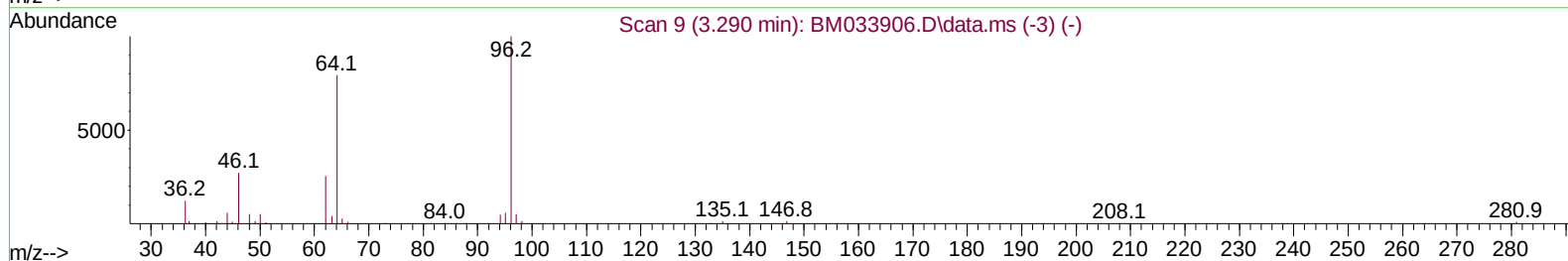
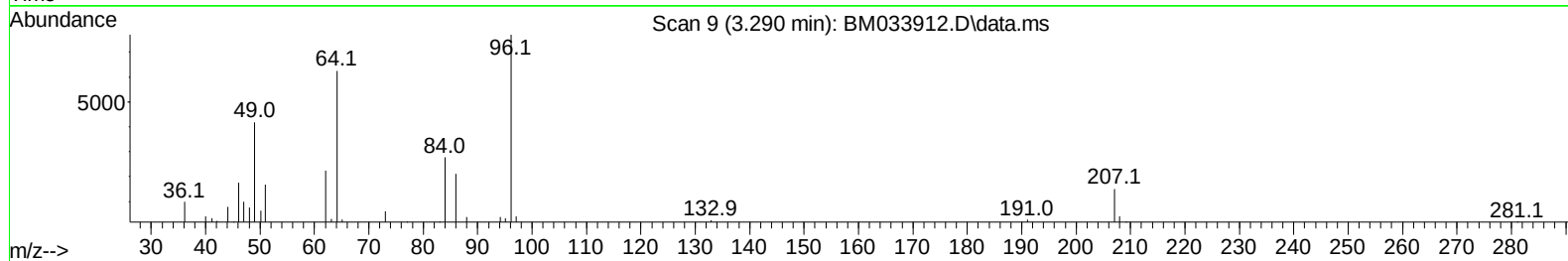
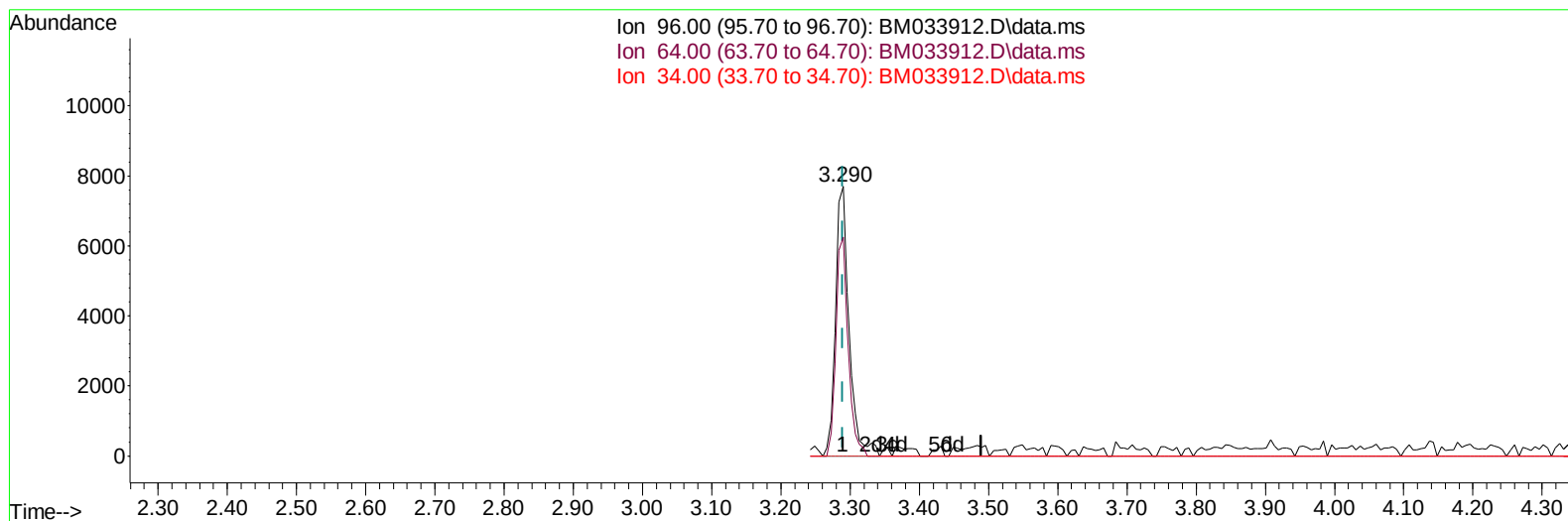
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TIC: BM033912.D\data.ms

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

3.290min (+ 0.000) 4.68 ng/uL m

response 10434

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	81.17
34.00	0.00	0.00
0.00	0.00	0.00

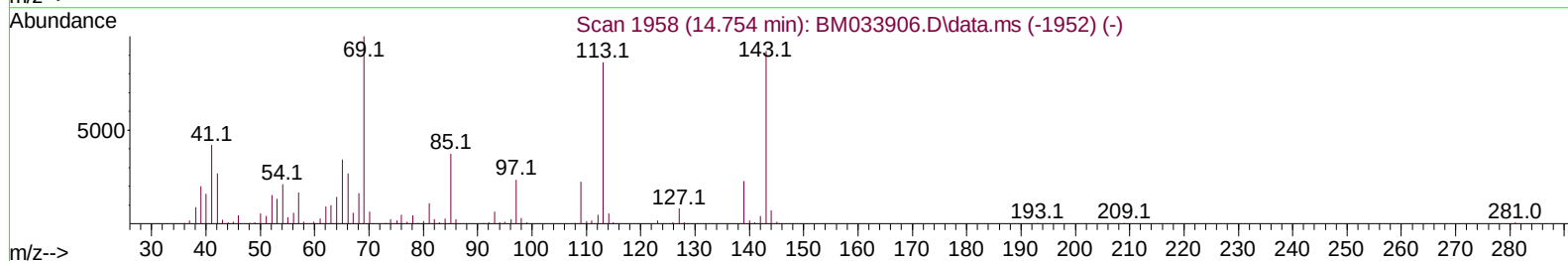
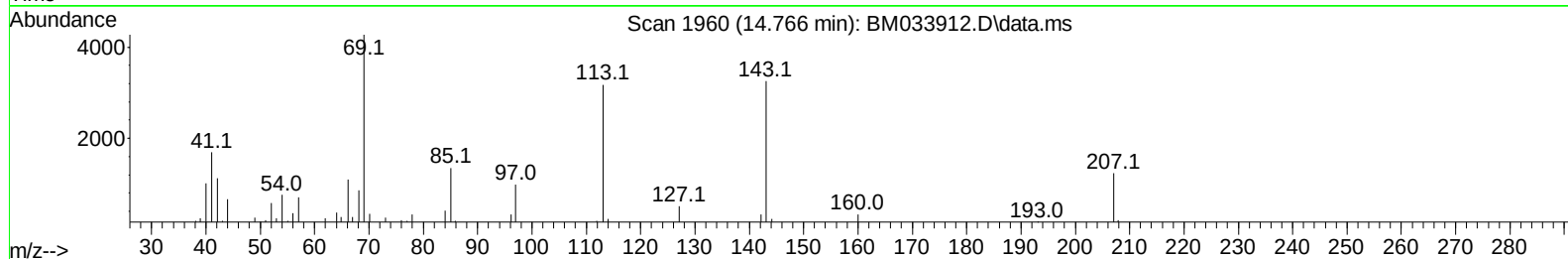
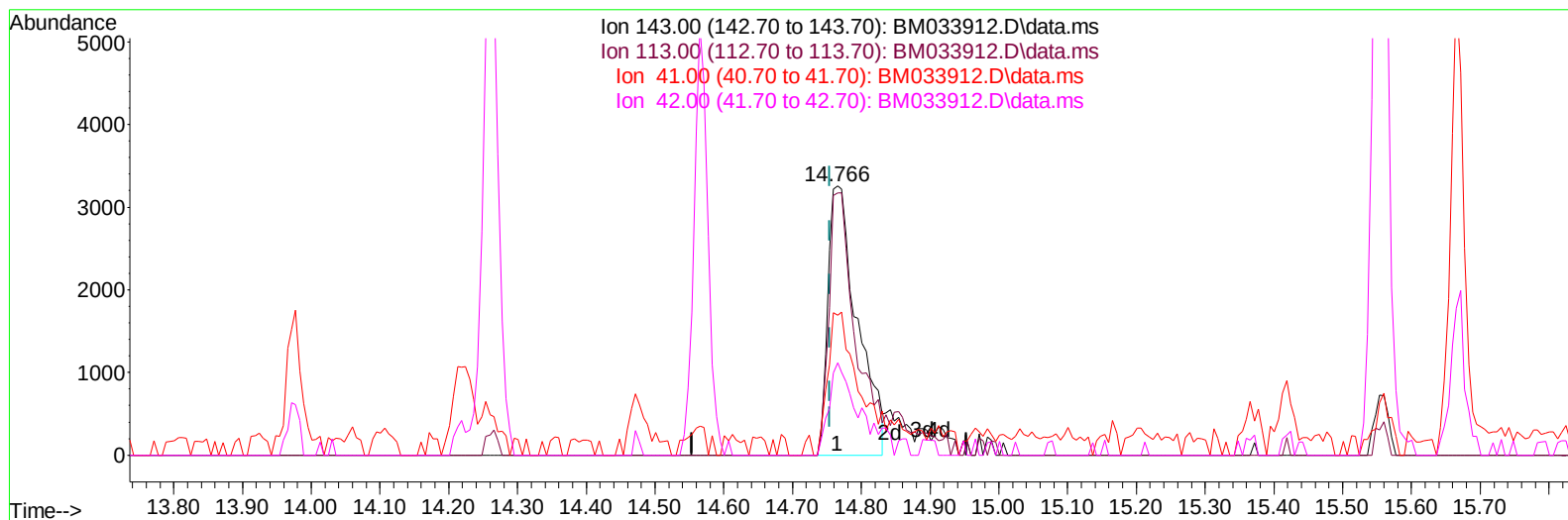
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0F28

Manual IntegrationsAPPROVED

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TIC: BM033912.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 2.97 ng/ul

response 9635

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	97.57
41.00	57.20	51.94
42.00	39.50	34.36

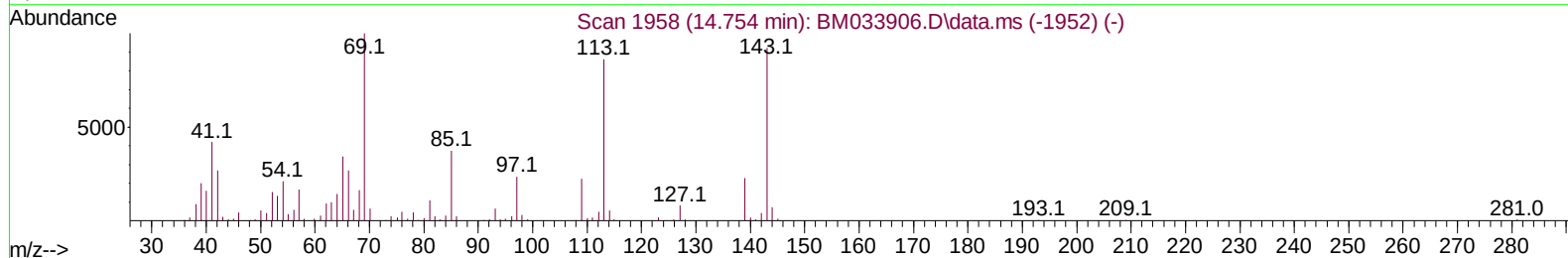
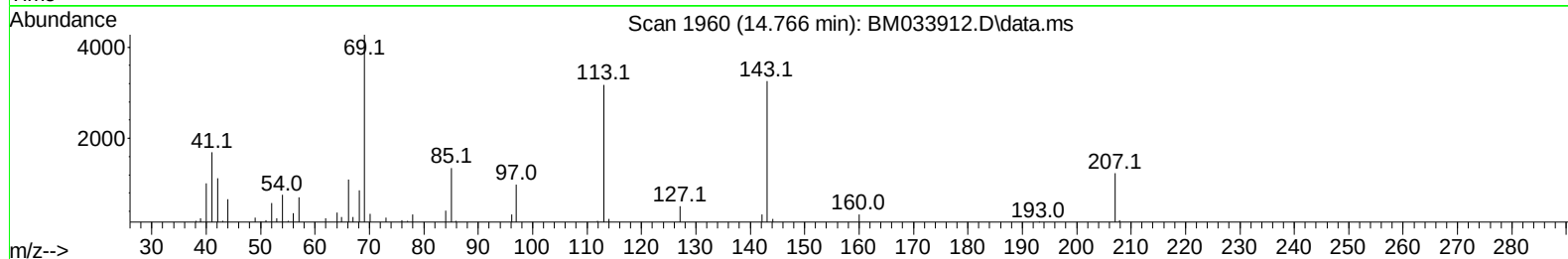
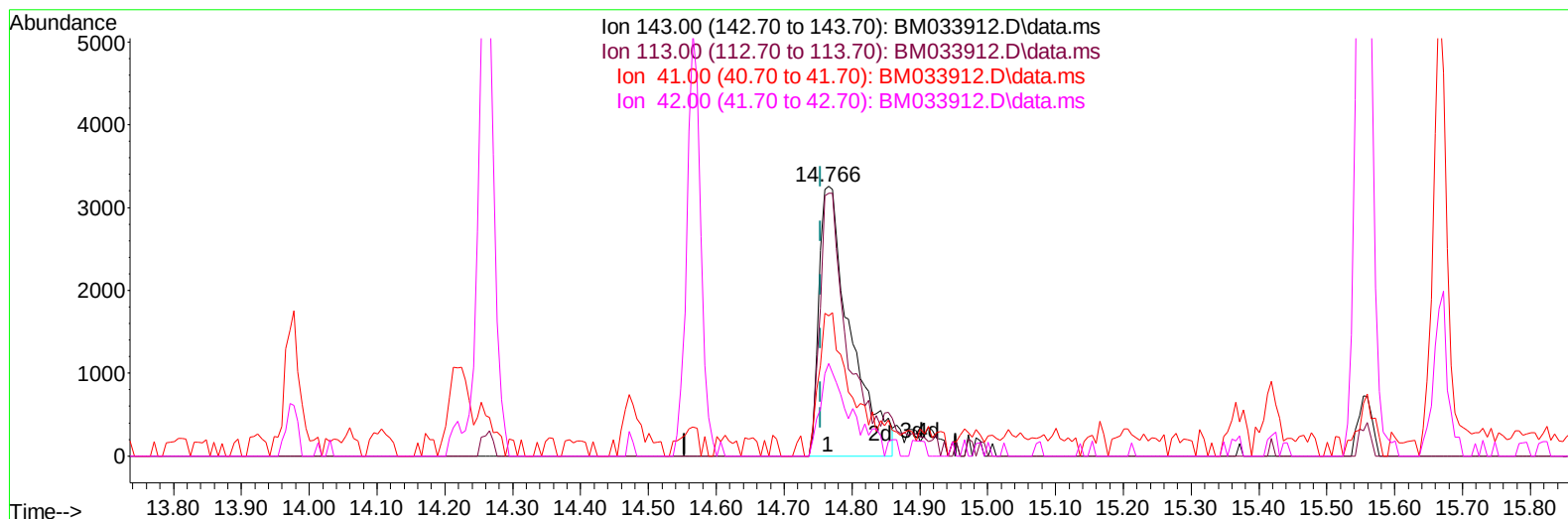
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Instrument :
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TIC: BM033912.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 3.22 ng/ul m

response 10435

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	97.57
41.00	57.20	51.94
42.00	39.50	34.36

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	83813	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	342741	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	214585	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	445500	20.000	ng/ul	0.00
79) Chrysene-d12	21.465	240	442012	20.000	ng/ul	0.00
88) Perylene-d12	23.859	264	443576	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	10434m	4.677	ng/uL	0.00
4) Pyridine-d5	3.719	84	53120	8.575	ng/ul	0.00
7) Phenol-d5	7.090	99	43553	5.786	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	160880	35.095	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	154689	26.878	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	88620	14.221	ng/ul	0.00
21) Nitrobenzene-d5	9.089	128	100349	37.074	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	101757	35.472	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	160538	29.192	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	233437	28.556	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	630141	37.272	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	731149	36.657	ng/ul	0.00
54) 4-Nitrophenol-d4	14.766	143	10435m	3.216	ng/ul	0.01
60) Fluorene-d10	15.554	176	536121	38.223	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	78027	29.444	ng/ul	0.00
73) Anthracene-d10	17.401	188	901870	41.720	ng/ul	0.00
81) Pyrene-d10	19.689	212	1028439	43.085	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.706	264	990611	42.921	ng/ul	0.00

Target Compounds Qvalue

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

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