

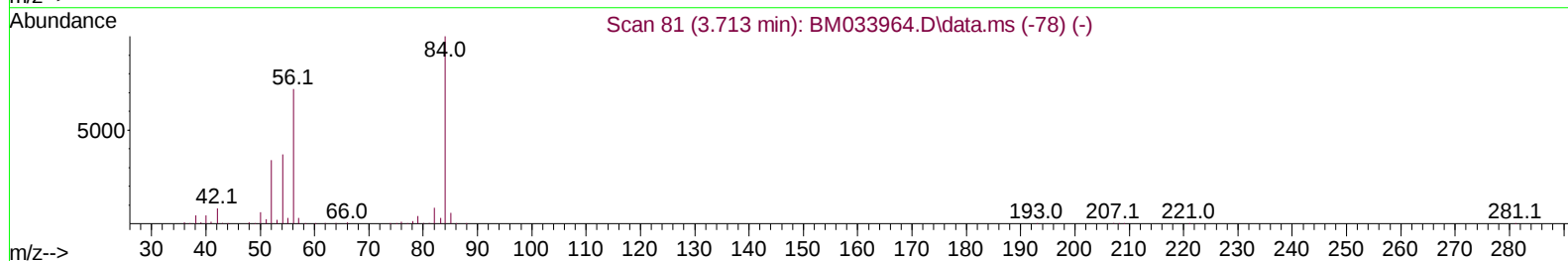
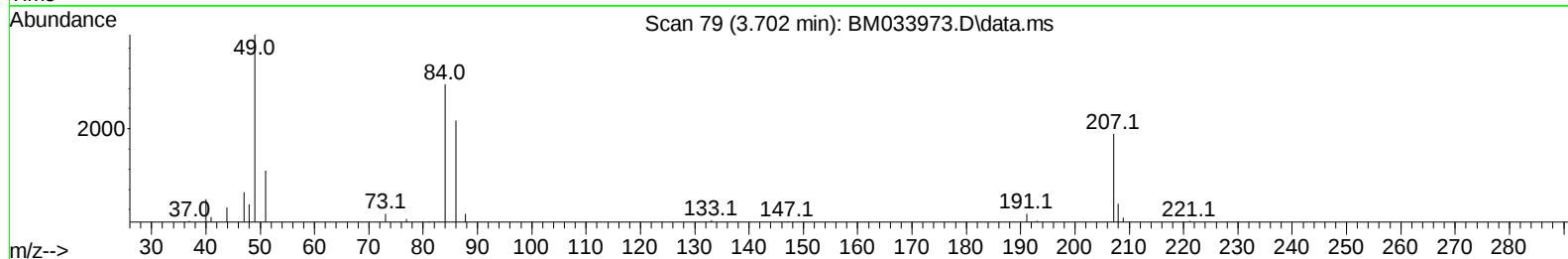
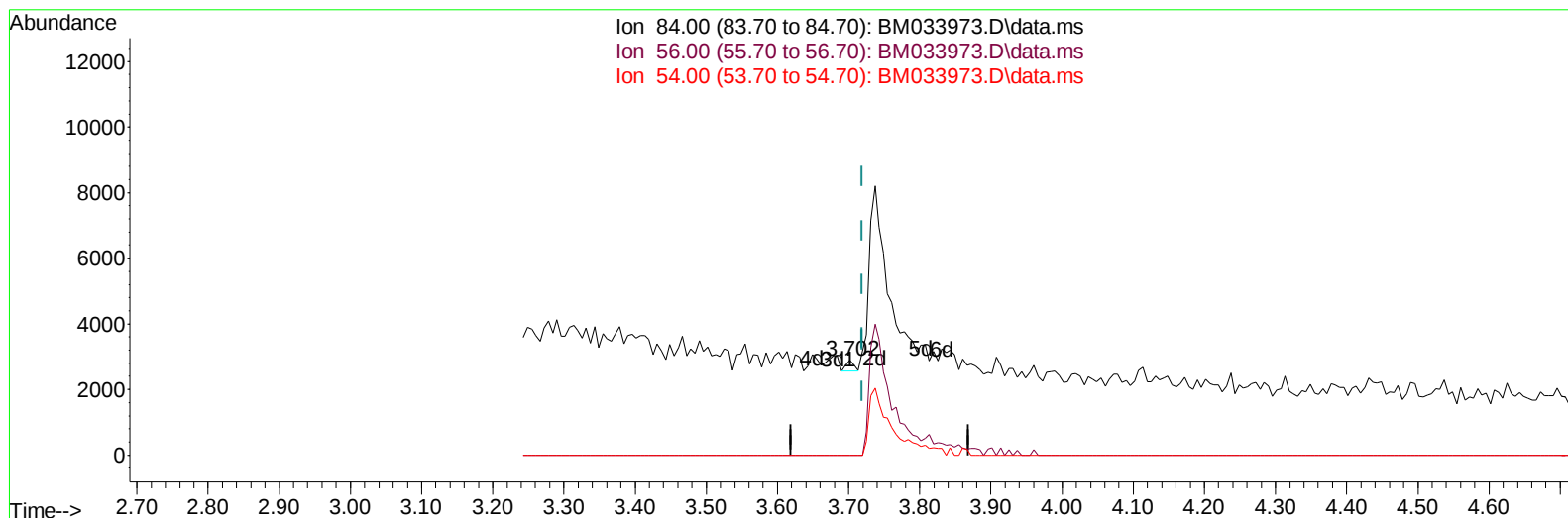
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033973.D
 Acq On : 02 Feb 2022 14:08
 Operator : CG/JU
 Sample : N1307-10DL 5X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q25DL

Manual IntegrationsAPPROVED

Quant Time: Feb 03 16:30:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 01 02:16:28 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/07/2022
 Supervised By :Yogesh Patel 02/08/2022



TIC: BM033973.D\data.ms

(4) Pyridine-d5 (S)

3.702min (-0.018) 0.03 ng/u1

response 250

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	0.00#
54.00	42.60	0.00#
0.00	0.00	0.00

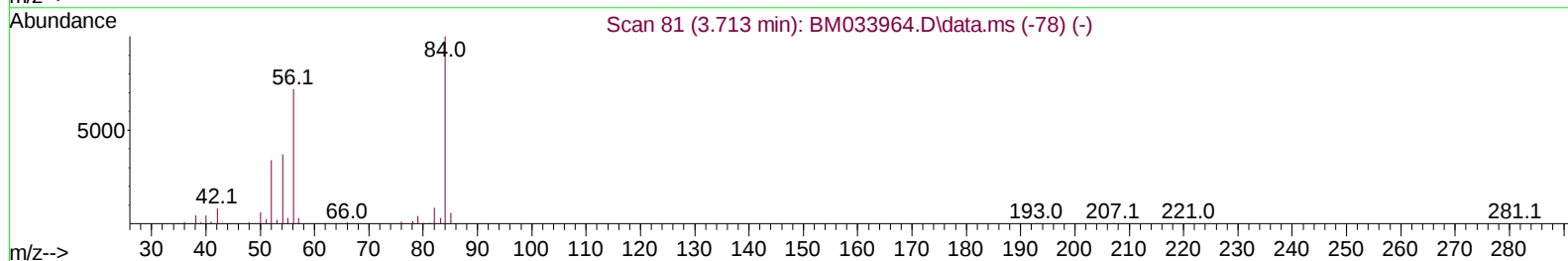
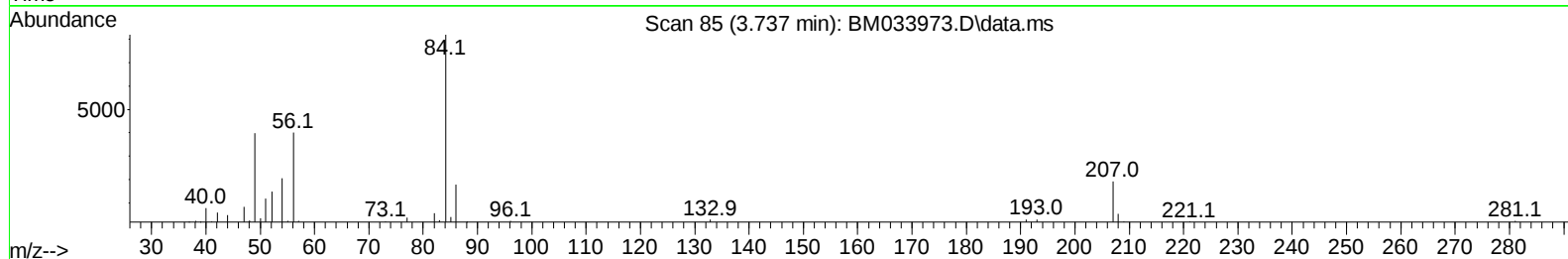
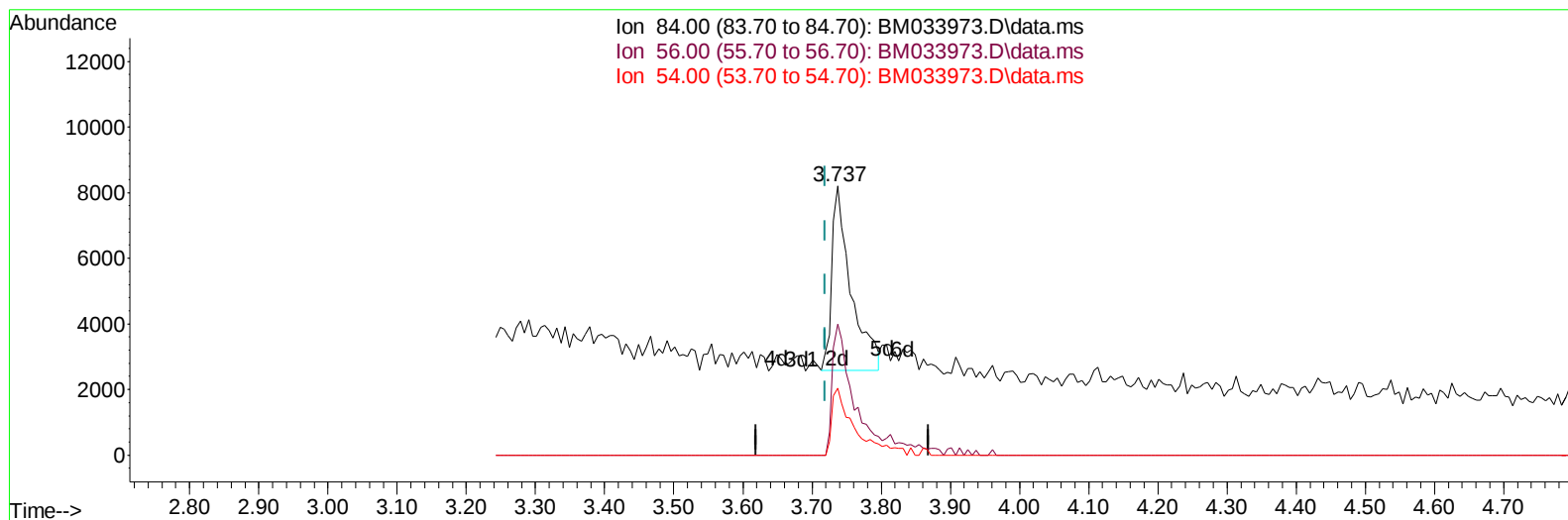
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(4) Pyridine-d5 (S)

3.737min (+ 0.018) 1.34 ng/ul m

response 10730

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	48.77#
54.00	42.60	25.00#
0.00	0.00	0.00

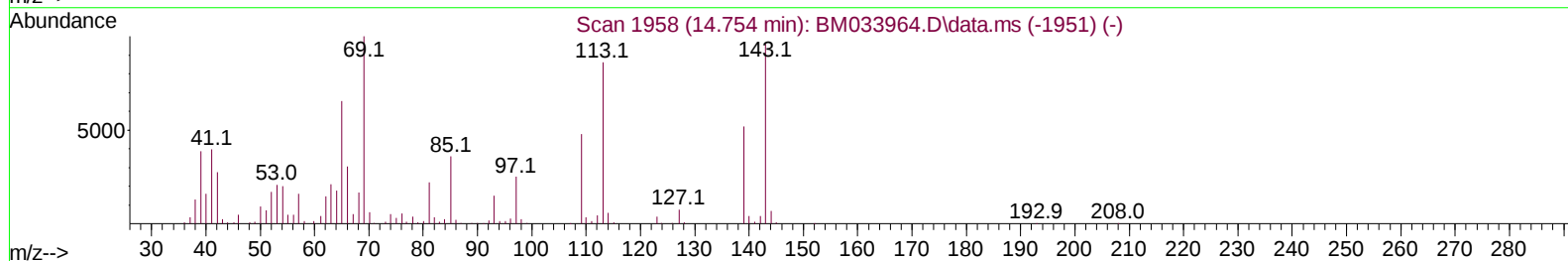
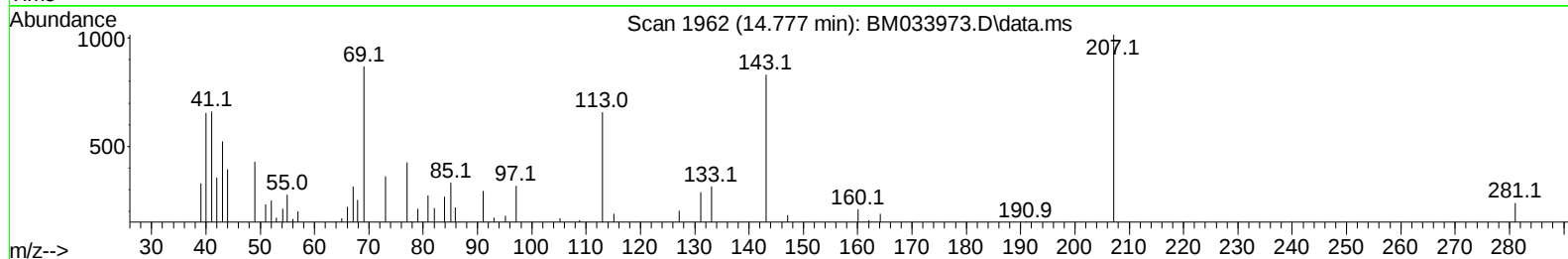
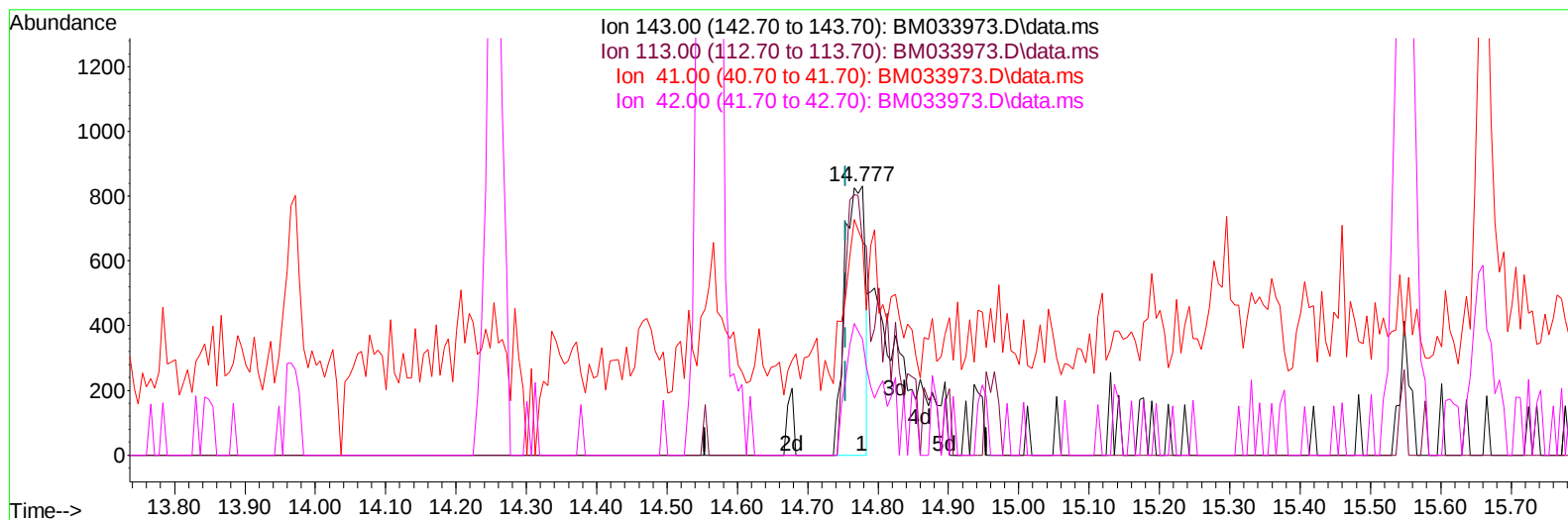
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TIC: BM033973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.777min (+ 0.024) 0.37 ng/u1

response 1694

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	79.18#
41.00	57.20	79.66#
42.00	39.50	42.96

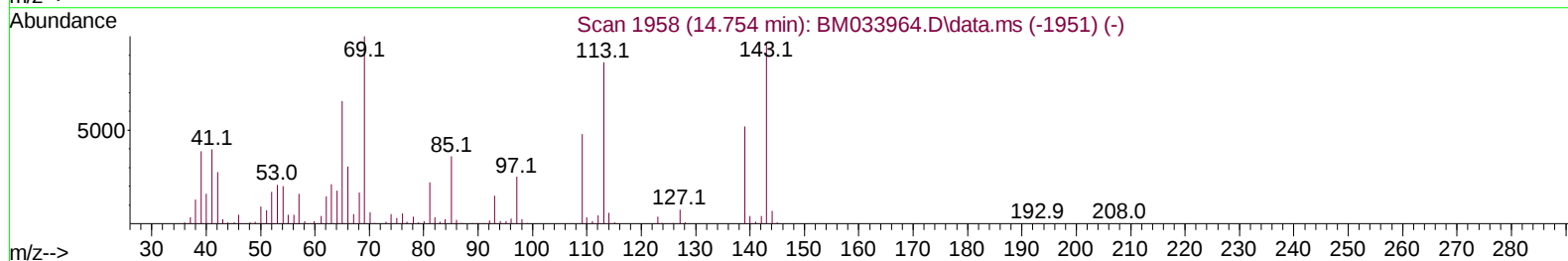
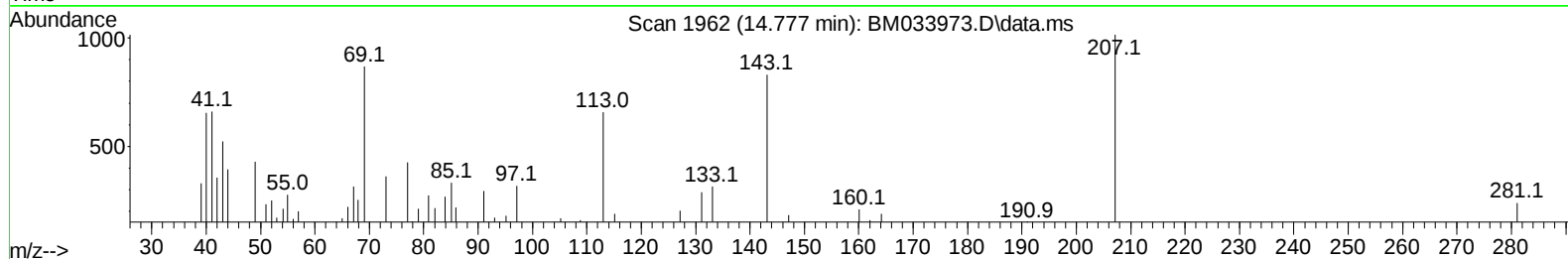
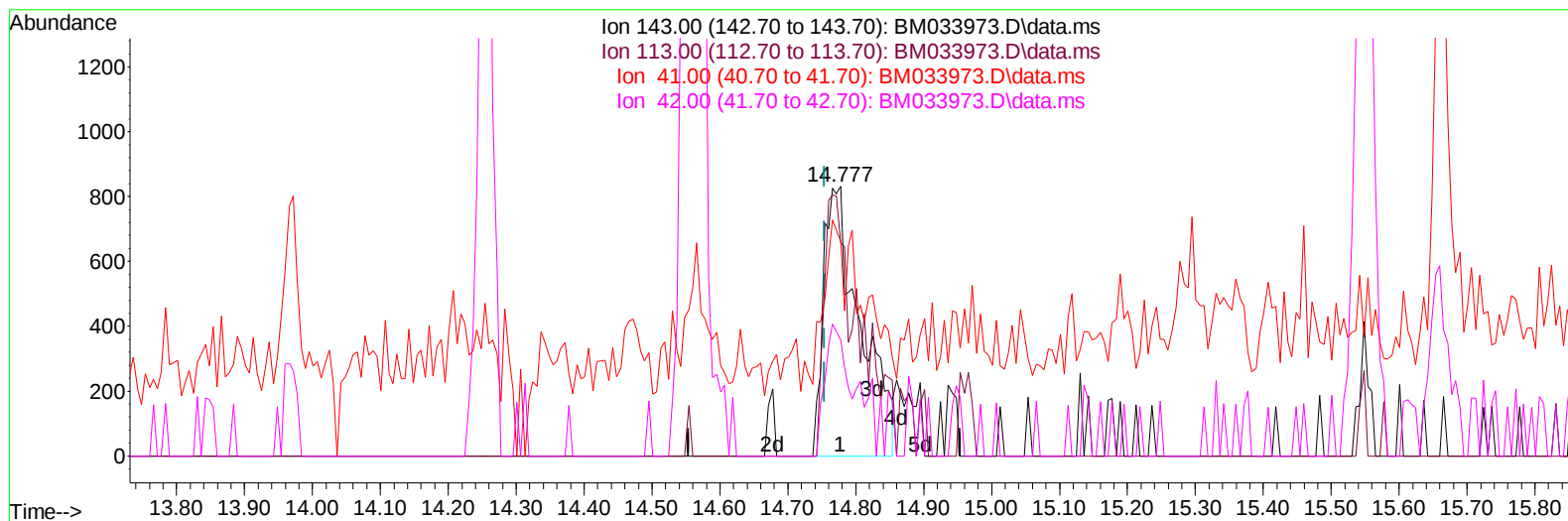
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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BM033973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.777min (+ 0.024) 0.68 ng/ul m

response 3122

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	79.18#
41.00	57.20	79.66#
42.00	39.50	42.96

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.925	152		112792	20.000	ng/ul	0.00
20) Naphthalene-d8	10.737	136		503080	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.560	164		359545	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188		799740	20.000	ng/ul	0.00
79) Chrysene-d12	21.459	240		748884	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264		624623	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.290	96		2344	0.820	ng/uL	0.00
4) Pyridine-d5	3.737	84		10730m	1.344	ng/ul	0.02
7) Phenol-d5	7.084	99		11489	1.216	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.249	67		37543	6.517	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132		35032	4.627	ng/ul	0.00
15) 4-Methylphenol-d8	8.631	113		22711	3.033	ng/ul	-0.01
21) Nitrobenzene-d5	9.084	128		21968	5.442	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143		21322	4.891	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165		39548	5.013	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131		44283	4.065	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166		201969	7.517	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160		211459	6.425	ng/ul	0.00
54) 4-Nitrophenol-d4	14.777	143		3122m	0.680	ng/ul	0.02
60) Fluorene-d10	15.548	176		168682	7.519	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.660	200		22050	4.392	ng/ul	-0.01
73) Anthracene-d10	17.395	188		283950	7.414	ng/ul	0.00
81) Pyrene-d10	19.677	212		338119	7.648	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.695	264		241310	7.355	ng/ul	-0.01
Target Compounds							
						Qvalue	
71) Pentachlorophenol	16.948	266		252008	39.767	ng/ul	97

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

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