

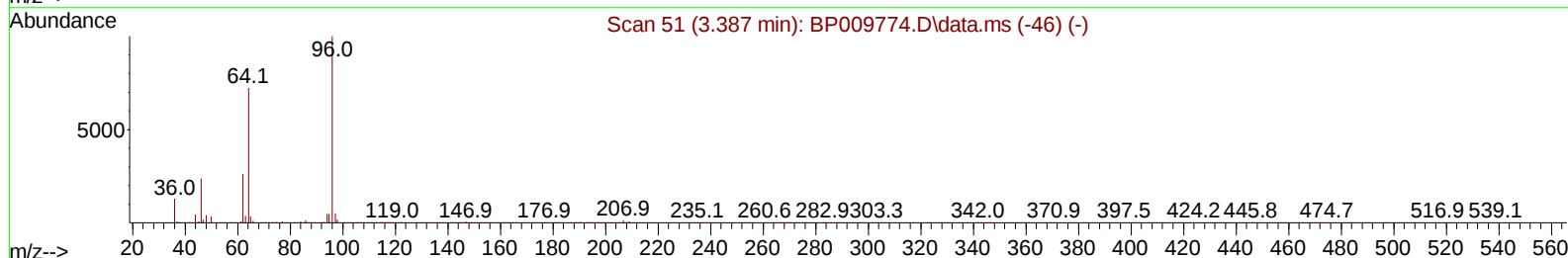
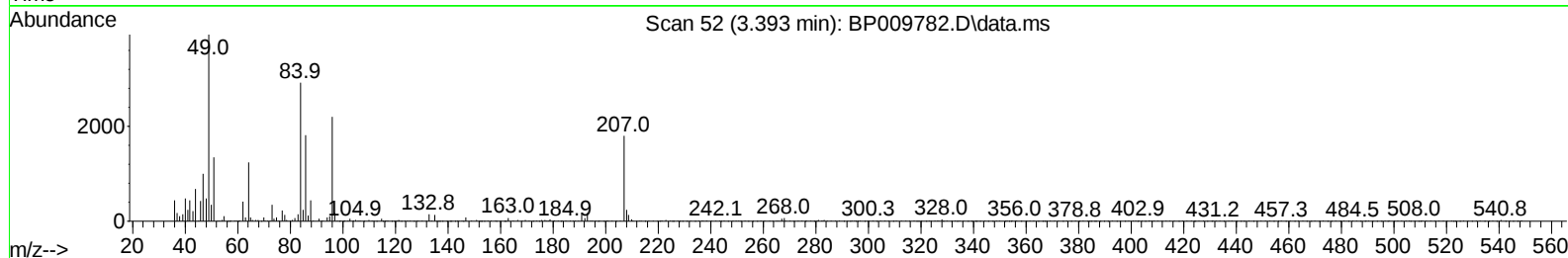
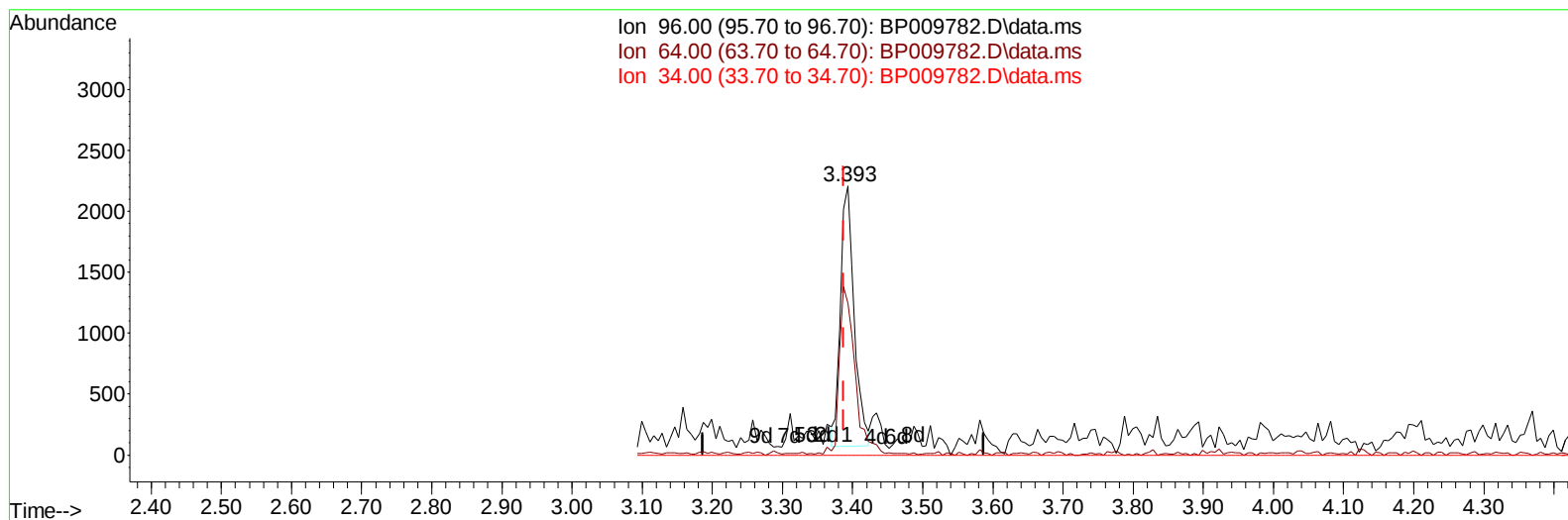
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009782.D
 Acq On : 12 Apr 2022 14:11
 Operator : CG/JU
 Sample : N2338-11DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J73DL

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:20:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009782.D\data.ms

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

3.393min (+ 0.006) 1.02 ng/uL

response 2908

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.39
34.00	0.00	0.00
0.00	0.00	0.00

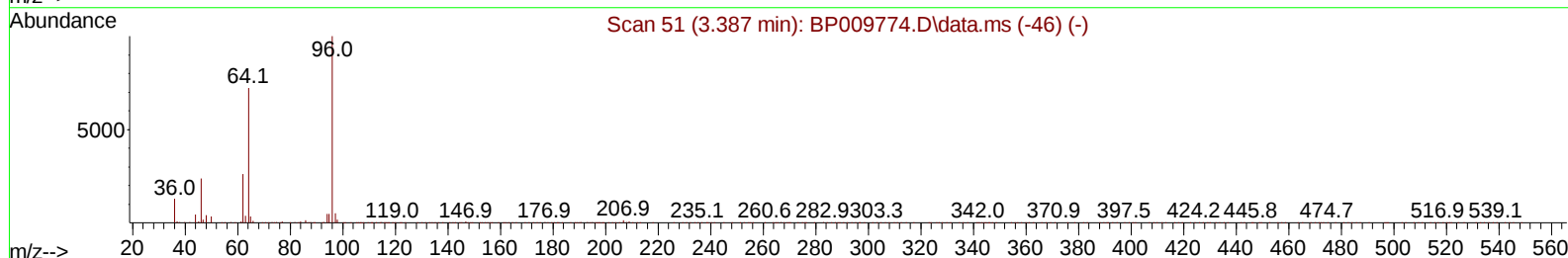
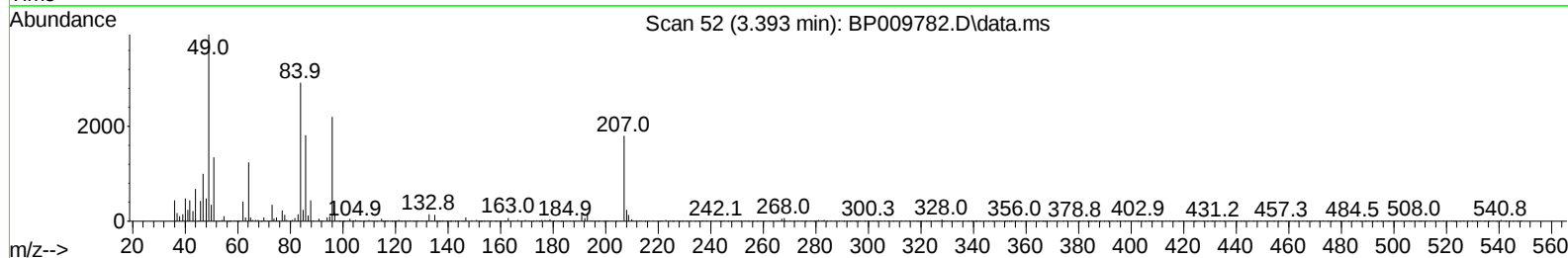
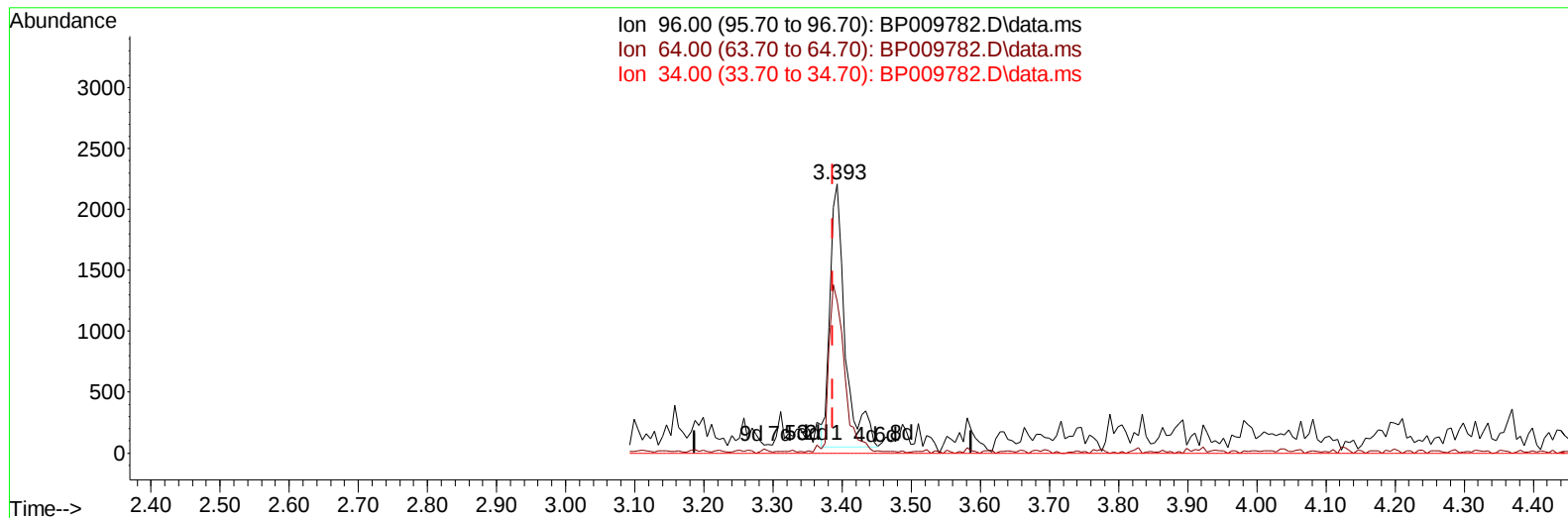
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TIC: BP009782.D\data.ms

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

3.393min (+ 0.006) 1.15 ng/uL m

response 3251

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.39
34.00	0.00	0.00
0.00	0.00	0.00

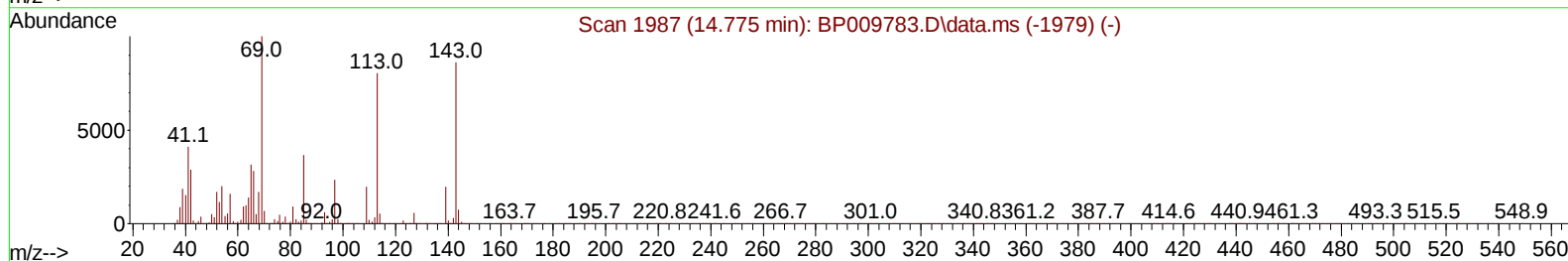
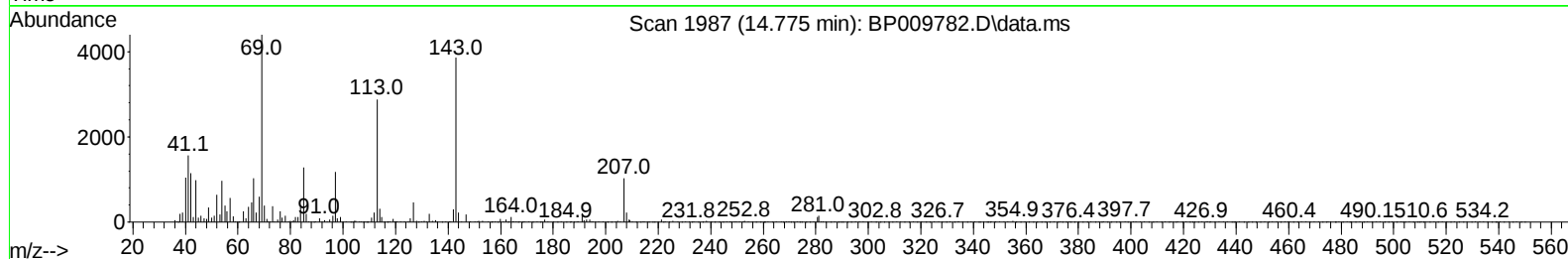
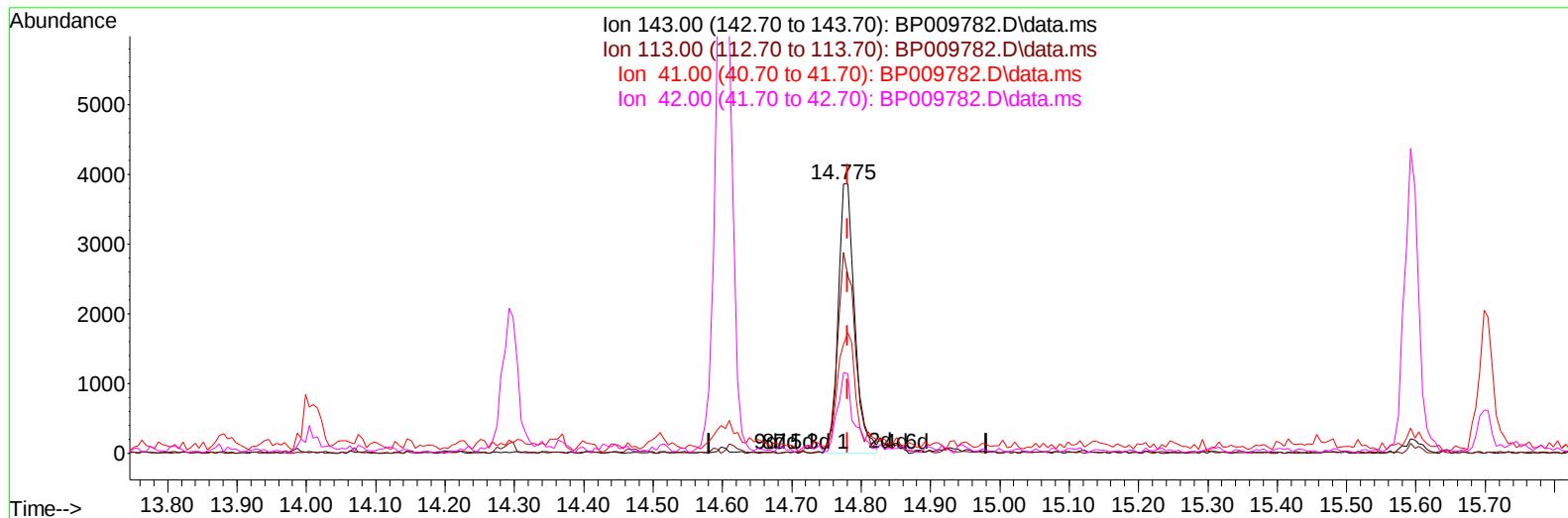
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TIC: BP009782.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 1.16 ng/u1

response 6307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.73#
41.00	23.20	40.44#
42.00	18.00	29.73#

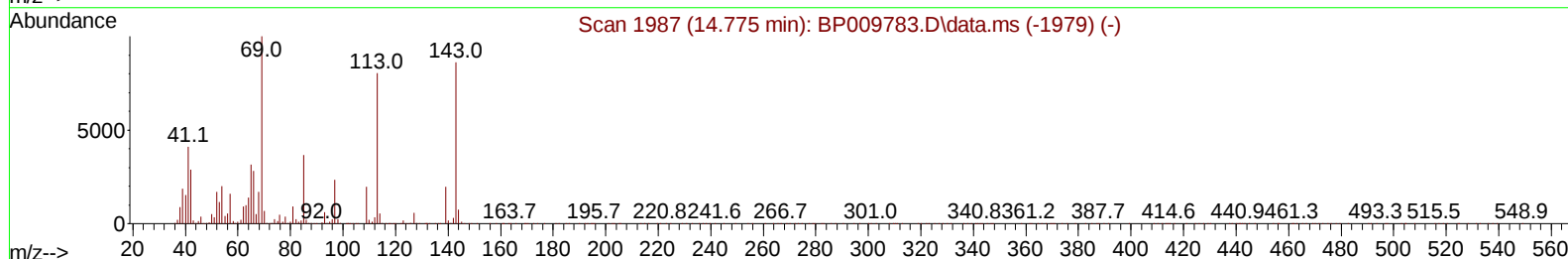
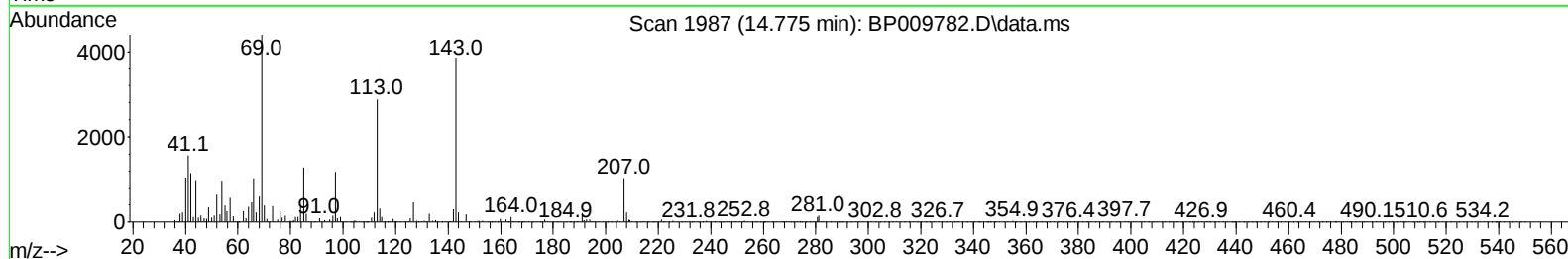
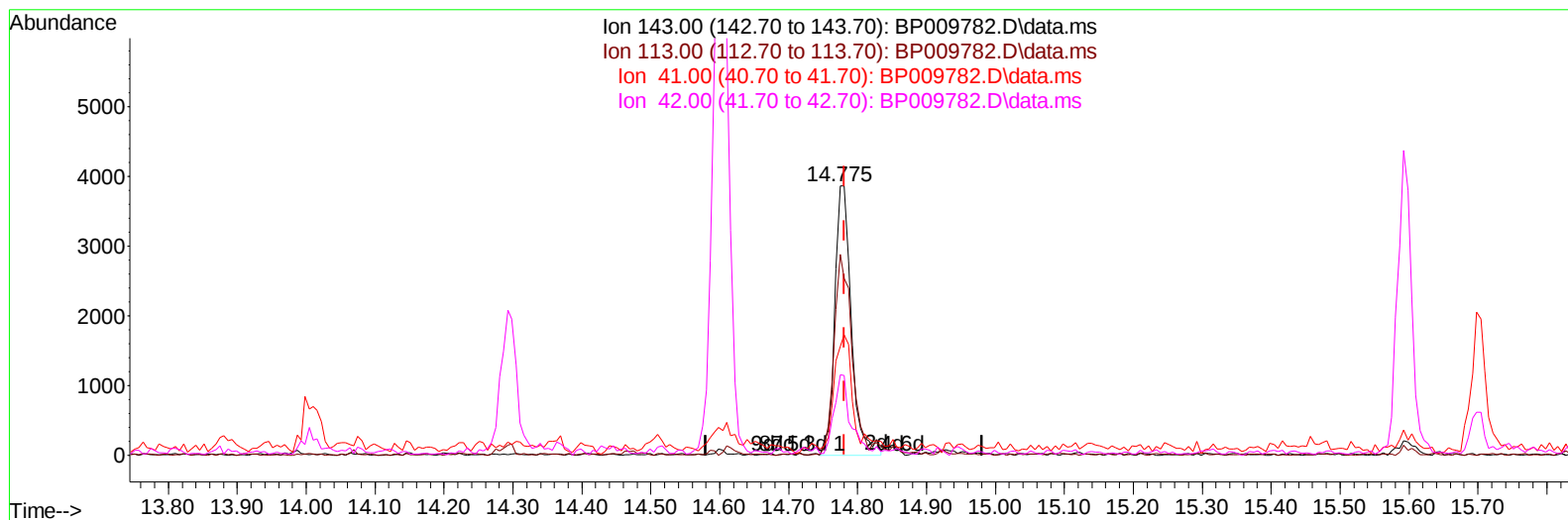
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 ALS Vial : 40 Sample Multiplier: 1

Instrument :
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TIC: BP009782.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 1.18 ng/ul m

response 6406

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.73#
41.00	23.20	40.44#
42.00	18.00	29.73#

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 Sample : N2338-11DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	126693	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	540519	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	381811	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	863149	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	960348	20.000	ng/ul	# 0.00
88) Perylene-d12	23.904	264	902470	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	3251m	1.145	ng/uL	0.00
4) Pyridine-d5	3.817	84	17400	2.223	ng/ul	0.01
7) Phenol-d5	7.122	99	15964	1.435	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	56804	8.578	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	47318	5.640	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	31061	3.345	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	32125	8.242	ng/ul	0.00
24) 2-Nitrophenol-d4	9.852	143	29615	6.930	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	53302	5.961	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	76287	5.990	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	230134	7.673	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	247005	6.952	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	6406m	1.178	ng/ul	0.00
60) Fluorene-d10	15.598	176	201860	8.001	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	26320	5.147	ng/ul	0.00
73) Anthracene-d10	17.451	188	347683	8.413	ng/ul	0.00
81) Pyrene-d10	19.680	212	433407	8.263	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.739	264	393737	8.378	ng/ul	0.00
Target Compounds						
					Qvalue	
22) Nitrobenzene	9.175	77	529378	52.912	ng/ul#	84
75) 1,2,3,4-Tetrachloroben...	13.404	216	13625	1.092	ng/uL#	81

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

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