

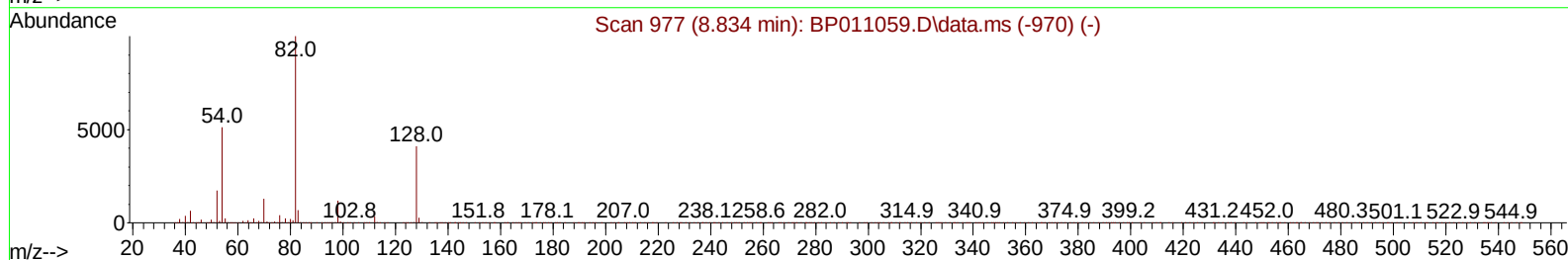
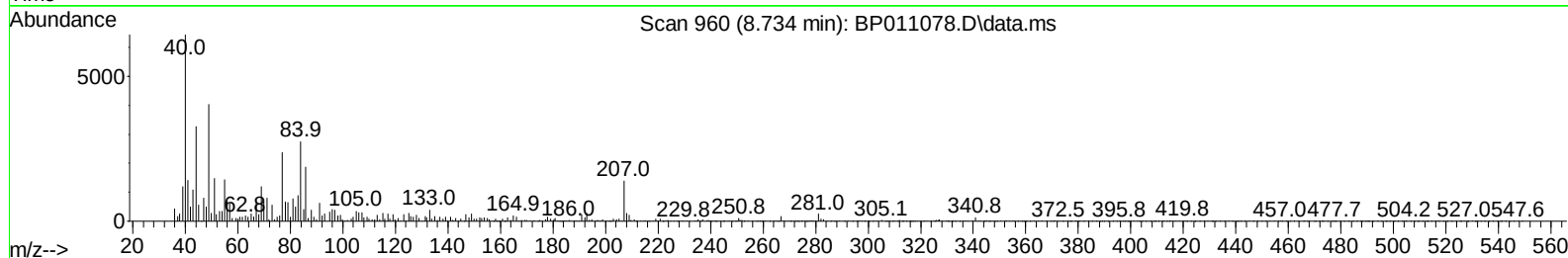
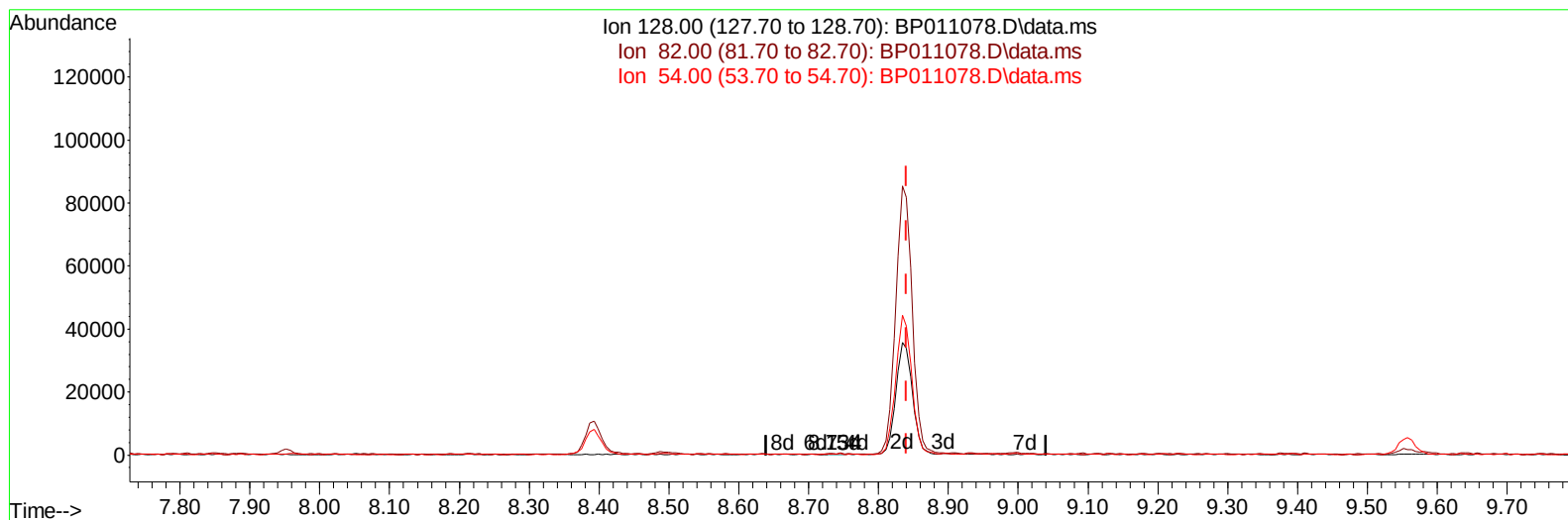
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011078.D
 Acq On : 22 Jul 2022 00:49
 Operator : CG/JU
 Sample : N3709-08 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0B24

Manual IntegrationsAPPROVED

Quant Time: Jul 22 01:35:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/22/2022
 Supervised By :mohammad ahmed 07/22/2022



TIC: BP011078.D\data.ms

(21) Nitrobenzene-d5 (S)

8.734min (-0.106) 0.00 ng/ul

response 77

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	233.64
54.00	155.90	159.35
0.00	0.00	0.00

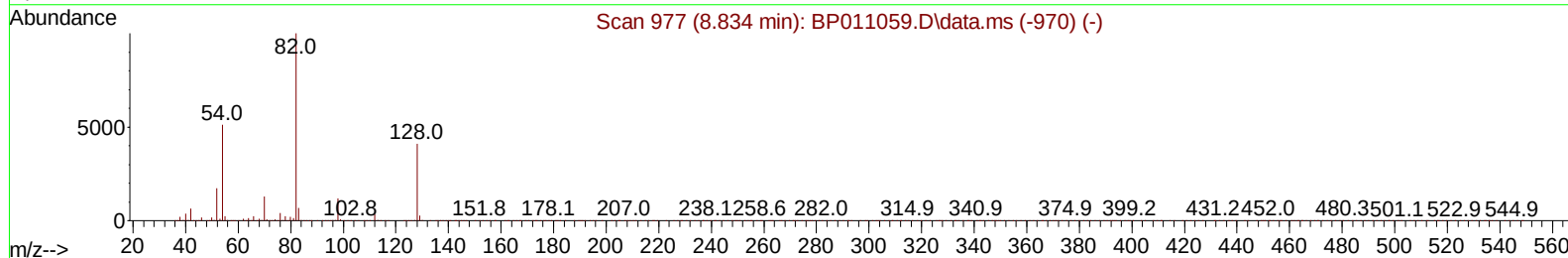
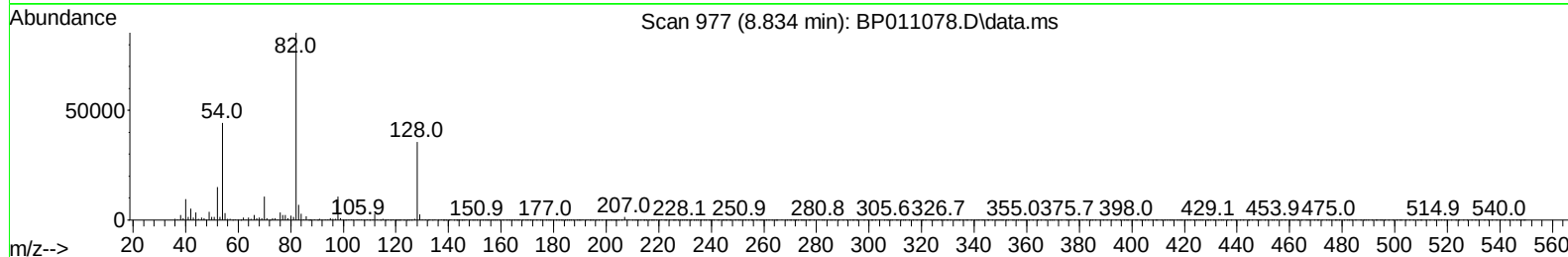
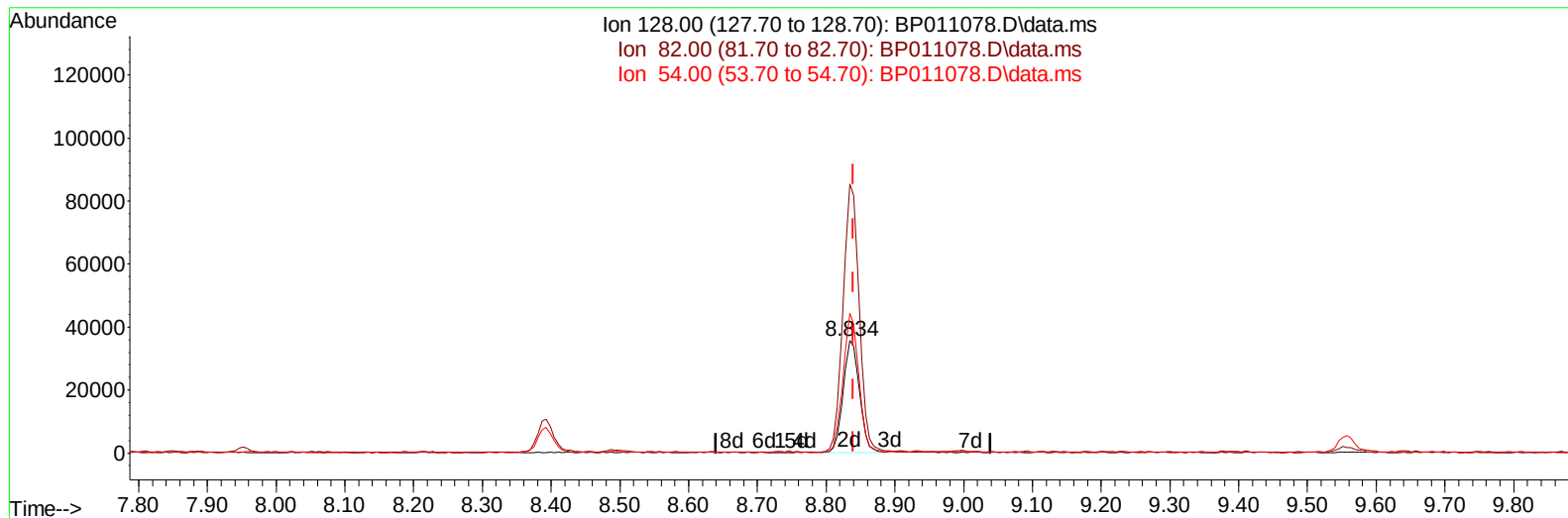
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(21) Nitrobenzene-d5 (S)

8.834min (-0.006) 3.28 ng/ul m

response 59060

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	238.94
54.00	155.90	124.03#
0.00	0.00	0.00

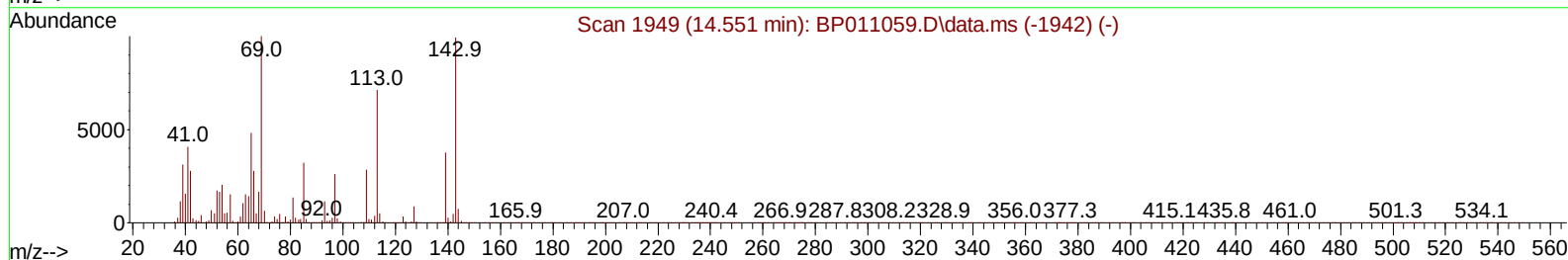
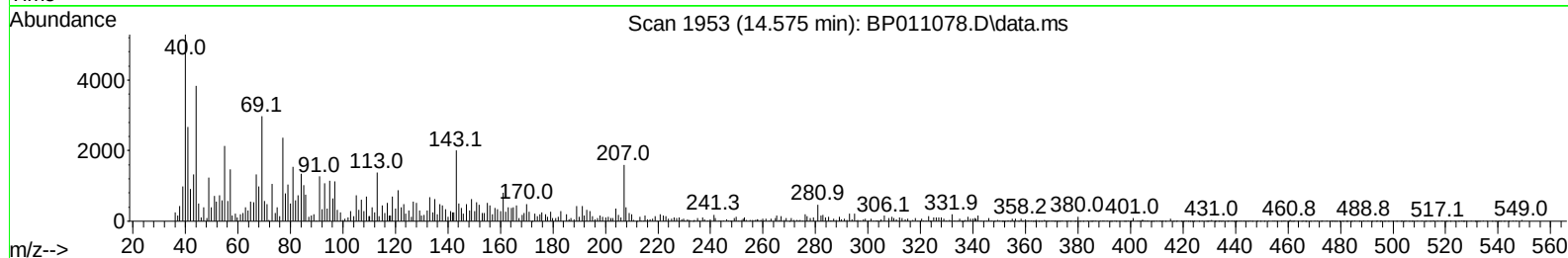
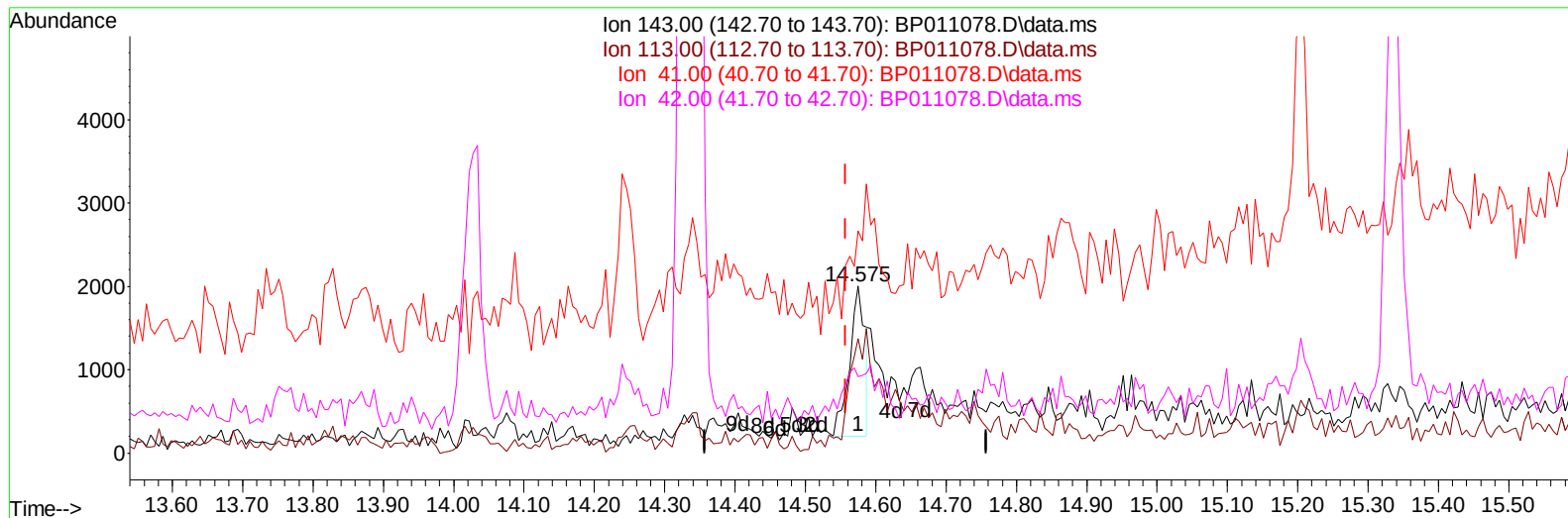
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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP011078.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.018) 0.14 ng/u1

response 2757

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	68.62#
41.00	59.00	133.28#
42.00	41.10	45.63

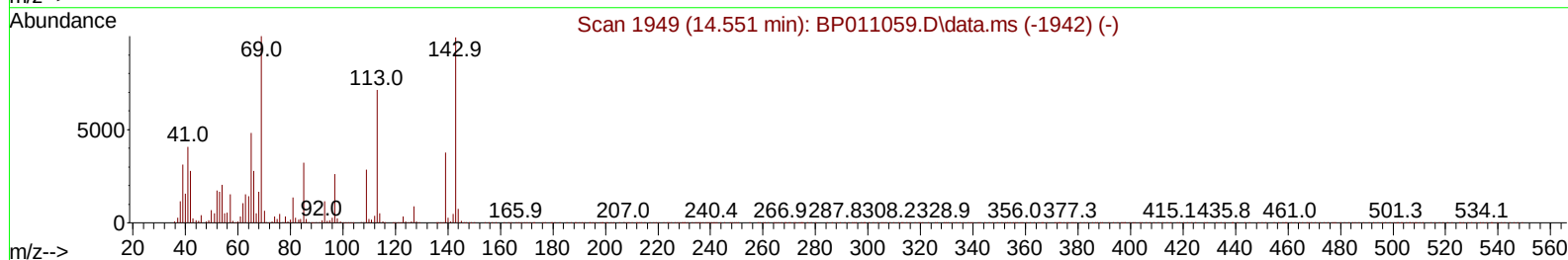
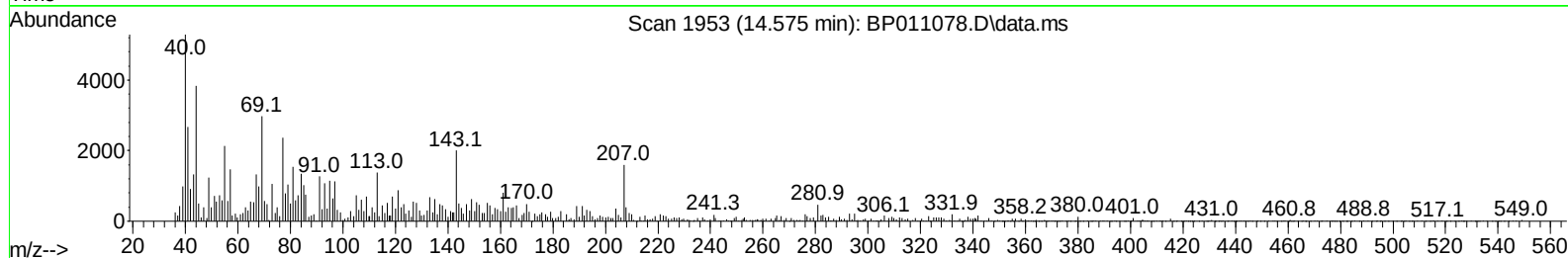
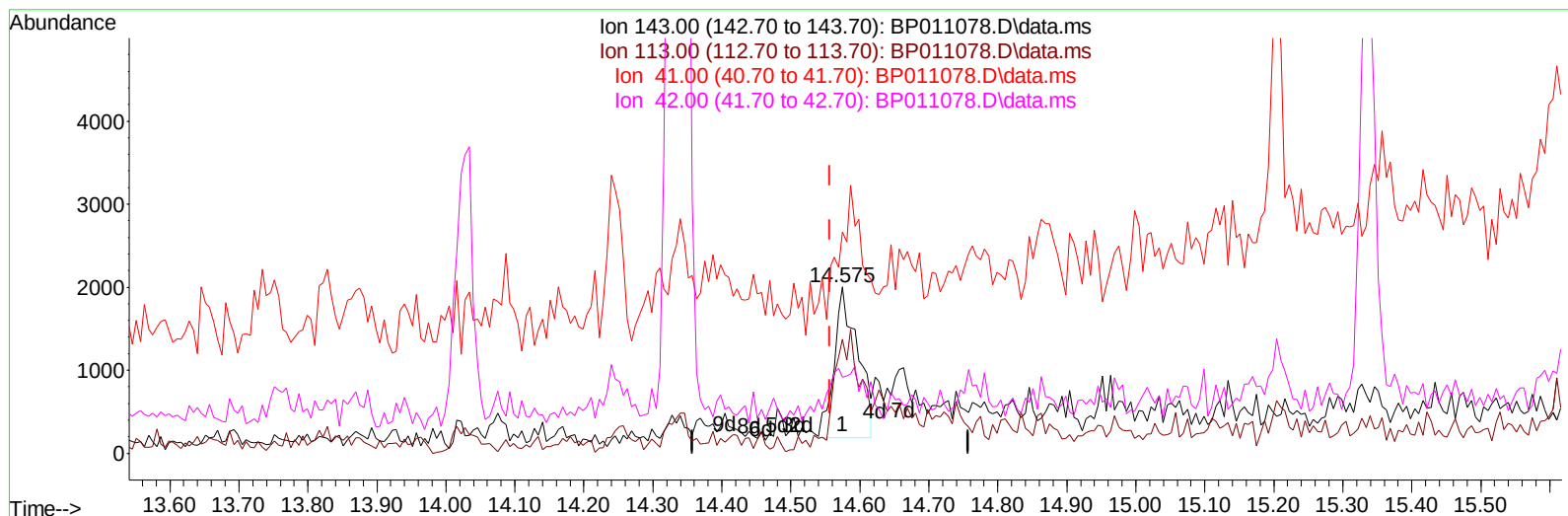
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TIC: BP011078.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.018) 0.22 ng/ul m

response 4347

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	68.62#
41.00	59.00	133.28#
42.00	41.10	45.63

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	590186	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	2474443	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.340	164	1417397	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	2616264	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1799414	20.000	ng/ul	# 0.00
88) Perylene-d12	23.586	264	1982335	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	3874	0.237	ng/uL	0.00
4) Pyridine-d5	3.581	84	770	0.018	ng/ul	-0.01
7) Phenol-d5	6.858	99	111881	2.309	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	99423	3.191	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	96708	2.470	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	93959	2.423	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	59060m	3.275	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	33378	1.871	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	73234	2.244	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	68232	1.268	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	247519	2.555	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160	439950	3.849	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	4347m	0.222	ng/ul	0.02
60) Fluorene-d10	15.340	176	316762	3.847	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	2522	0.195	ng/ul	0.00
73) Anthracene-d10	17.204	188	449900	4.045	ng/ul	0.00
81) Pyrene-d10	19.463	212	420264	4.447	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.427	264	374074	3.929	ng/ul	0.00

Target Compounds Qvalue

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

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