

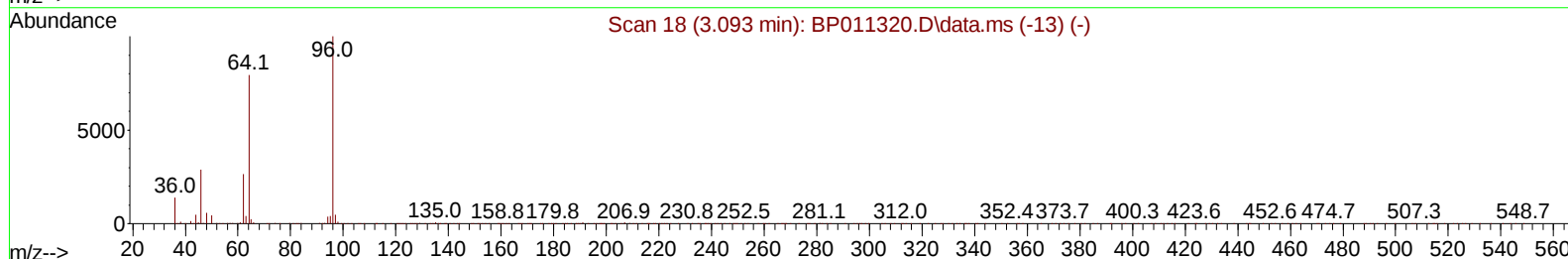
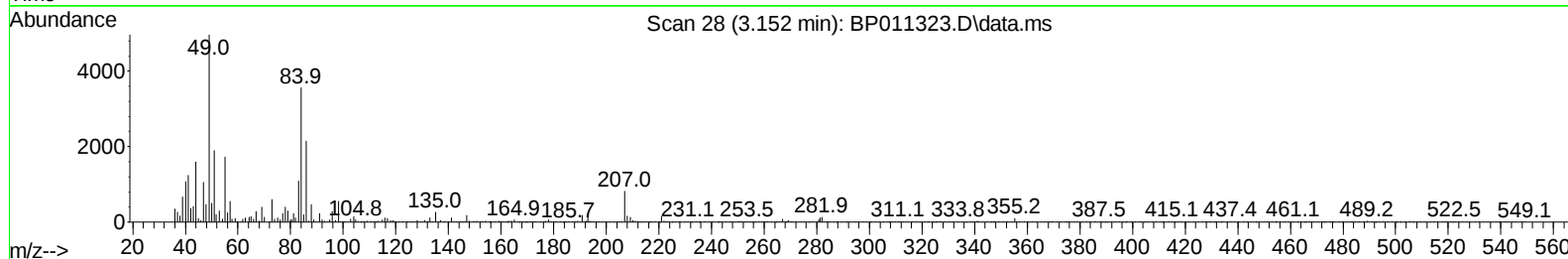
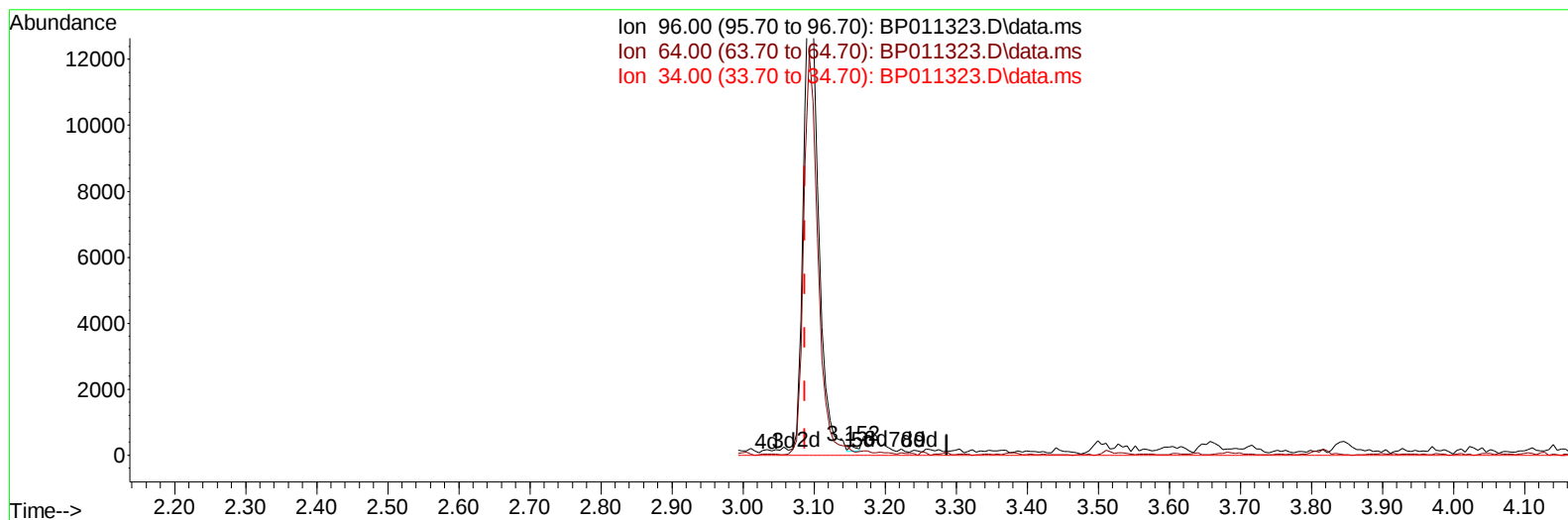
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011323.D
 Acq On : 09 Aug 2022 11:26
 Operator : CG/JU
 Sample : N3900-11
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C88

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:09:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 18:47:19 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/10/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011323.D\data.ms

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

3.152min (+ 0.065) 0.03 ng/uL

response	121	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	49.82
34.00	0.00	0.00
0.00	0.00	0.00

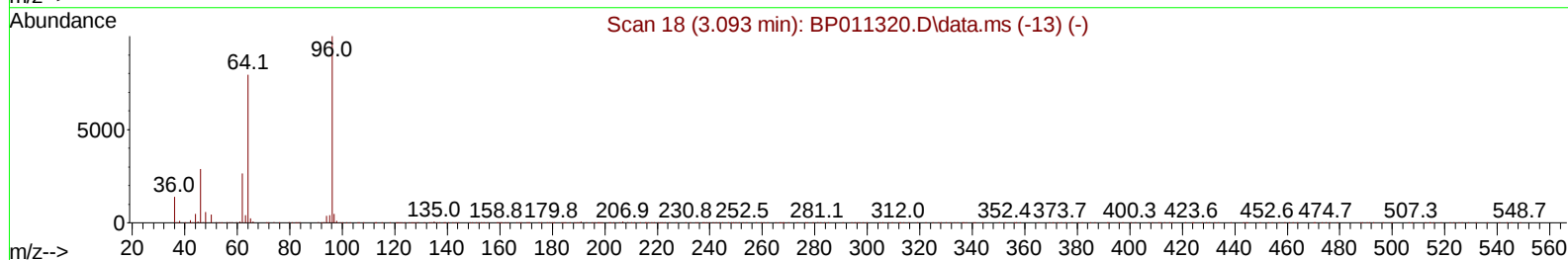
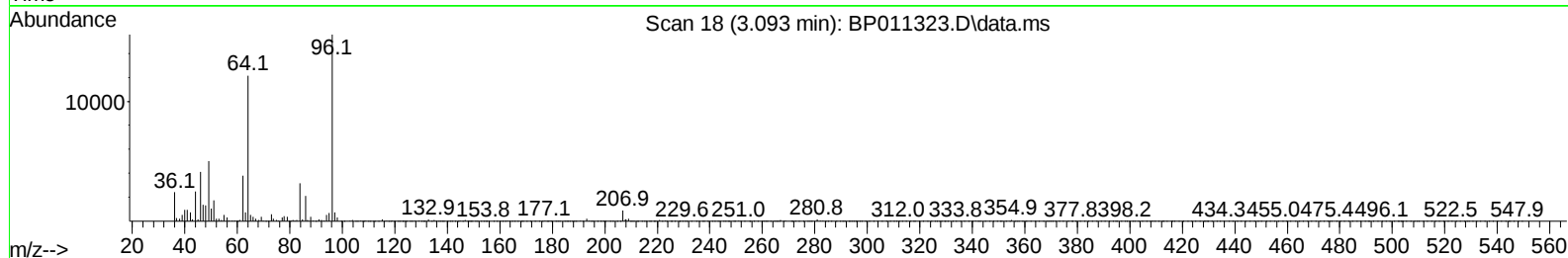
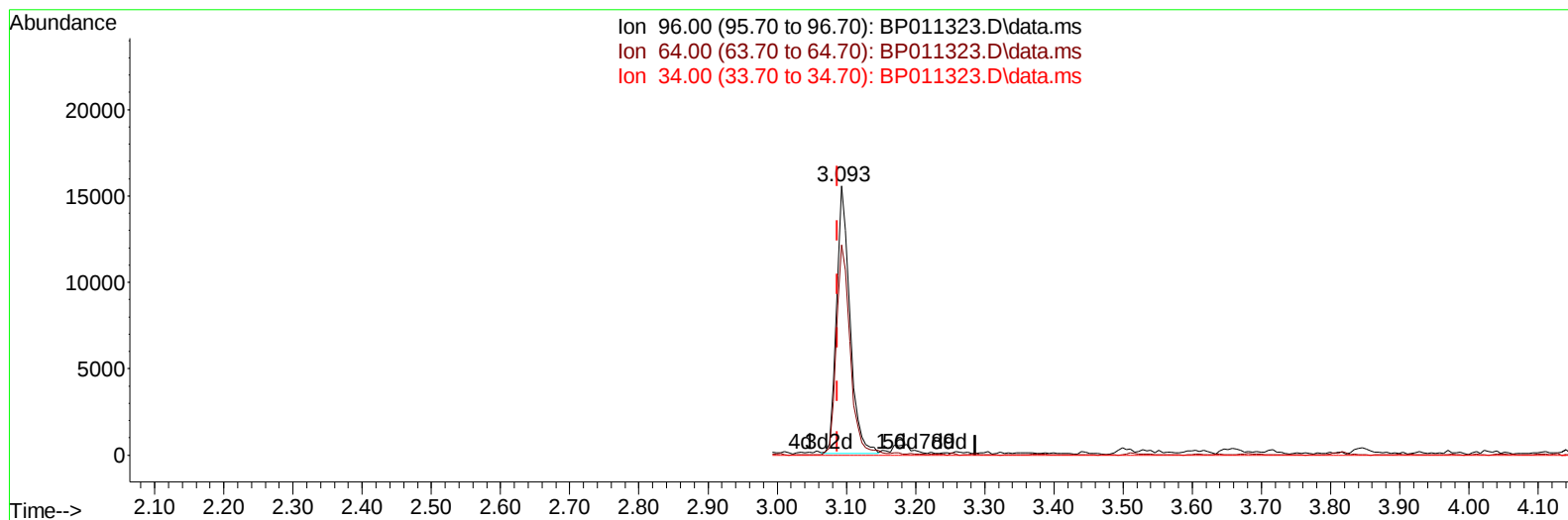
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TIC: BP011323.D\data.ms

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

3.093min (+ 0.006) 4.42 ng/uL m

response 20905

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	78.09#
34.00	0.00	0.00
0.00	0.00	0.00

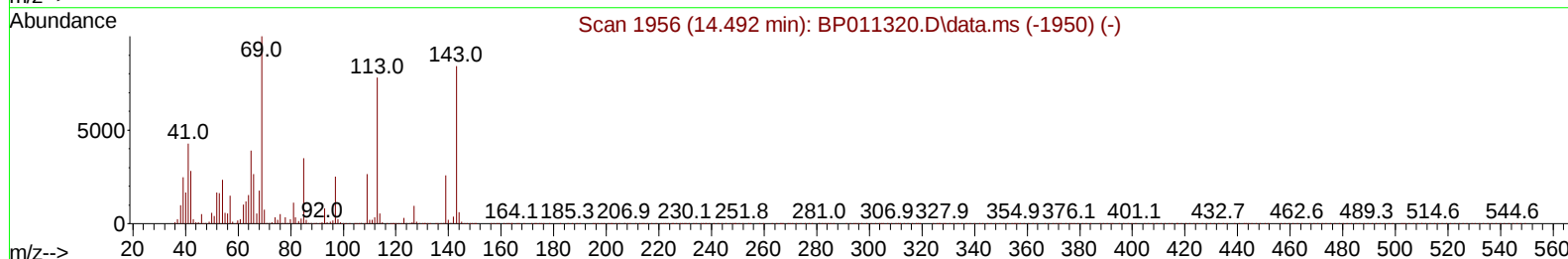
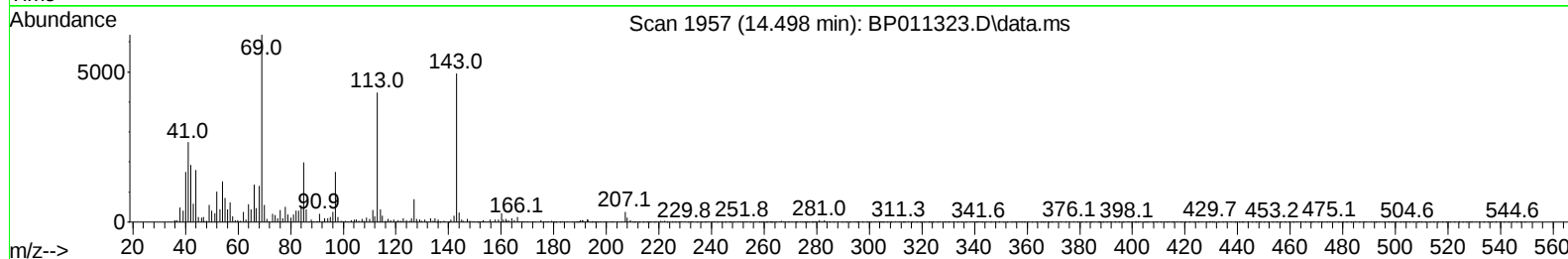
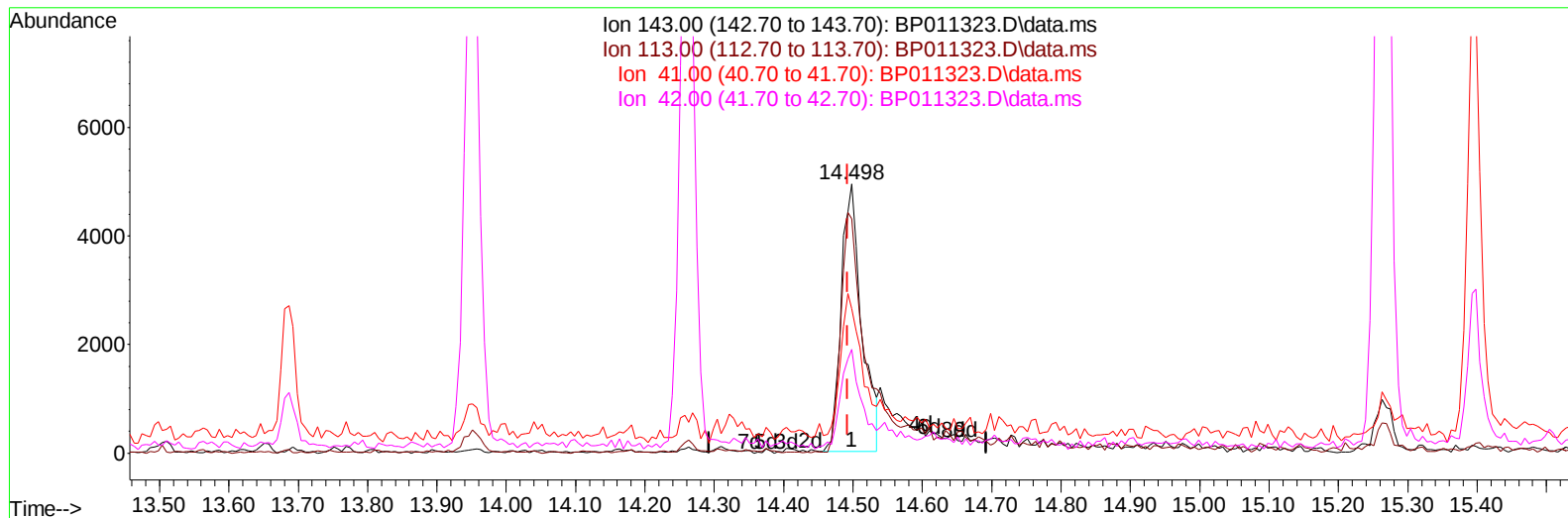
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TIC: BP011323.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.24 ng/u1

response 9981

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	87.18
41.00	54.40	53.77
42.00	35.50	38.50

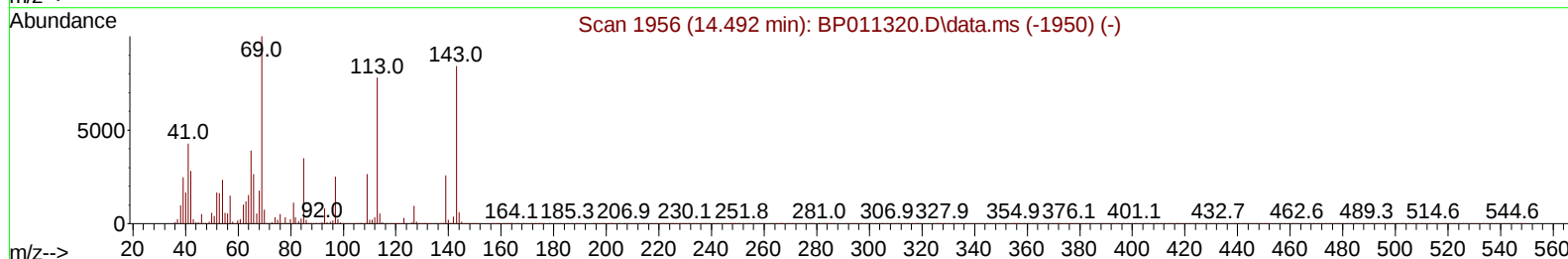
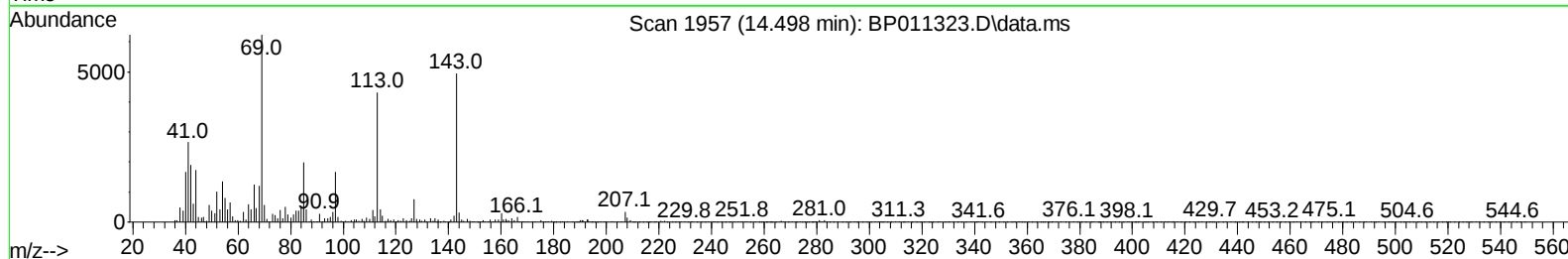
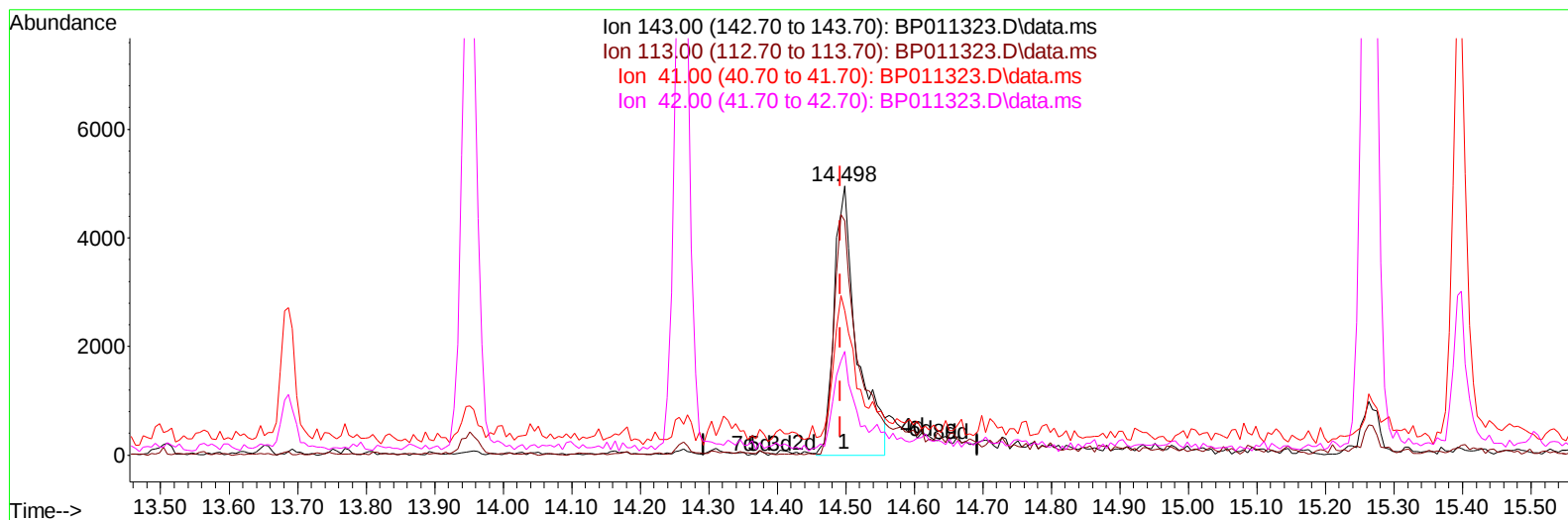
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Instrument :
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TIC: BP011323.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.56 ng/ul m

response 11419

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	87.18
41.00	54.40	53.77
42.00	35.50	38.50

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	145050	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	615507	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	351881	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	705850	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	661523	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	666403	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	20905m	4.418	ng/uL	0.00
4) Pyridine-d5	3.517	84	45872	3.843	ng/ul	0.02
7) Phenol-d5	6.775	99	57183	4.455	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	234950	25.814	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	188763	18.742	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	113840	11.295	ng/ul	0.00
21) Nitrobenzene-d5	8.757	128	123867	29.570	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	103182	26.904	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	179037	23.116	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	228732	16.893	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	775170	31.749	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	861888	29.326	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	11419m	2.561	ng/ul	0.00
60) Fluorene-d10	15.263	176	640556	30.678	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	85642	26.122	ng/ul	0.00
73) Anthracene-d10	17.133	188	1138738	35.892	ng/ul	0.00
81) Pyrene-d10	19.392	212	1315666	36.099	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1222544	36.378	ng/ul	0.00

Target Compounds Qvalue

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

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