

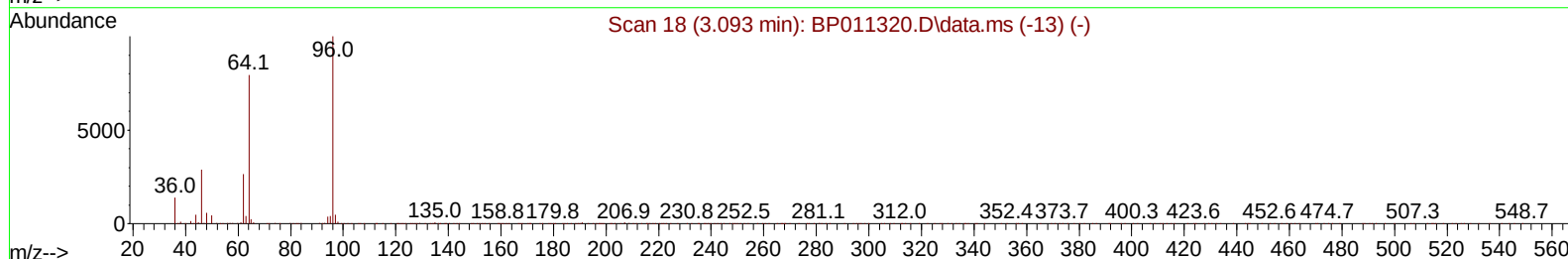
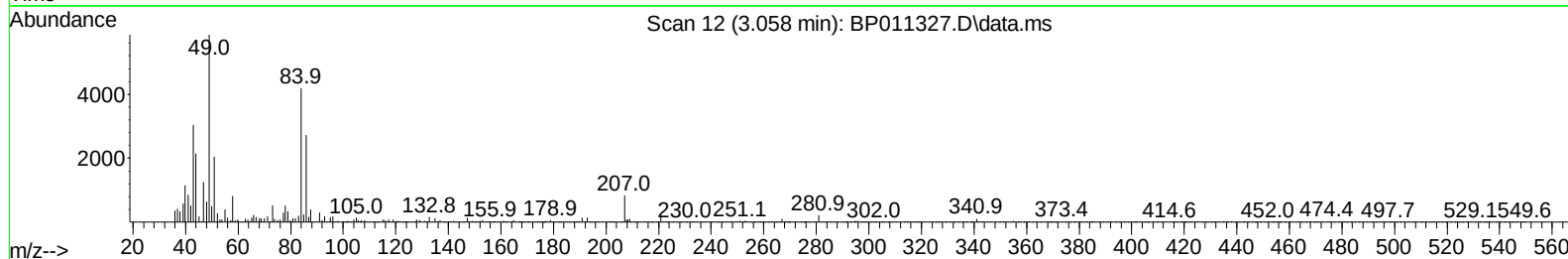
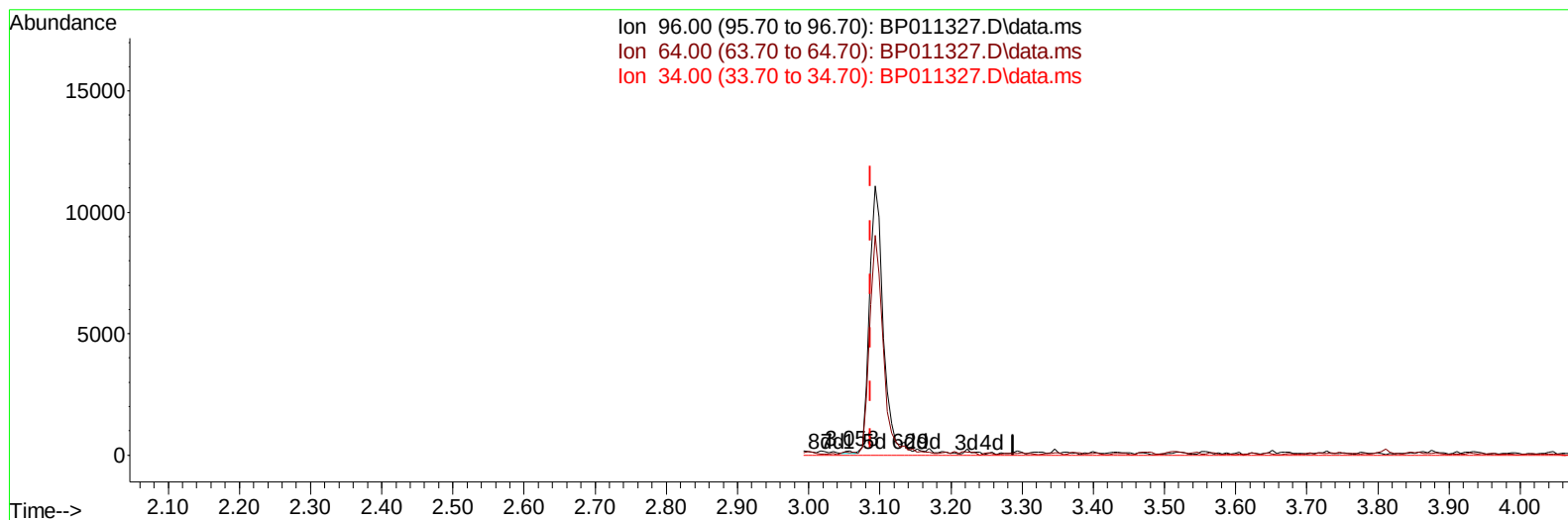
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011327.D
 Acq On : 09 Aug 2022 13:45
 Operator : CG/JU
 Sample : N4026-23
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJJ8

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:10:53 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: BP011327.D\data.ms

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

3.058min (-0.029) 0.02 ng/uL

response	85	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	58.72
34.00	0.00	0.00
0.00	0.00	0.00

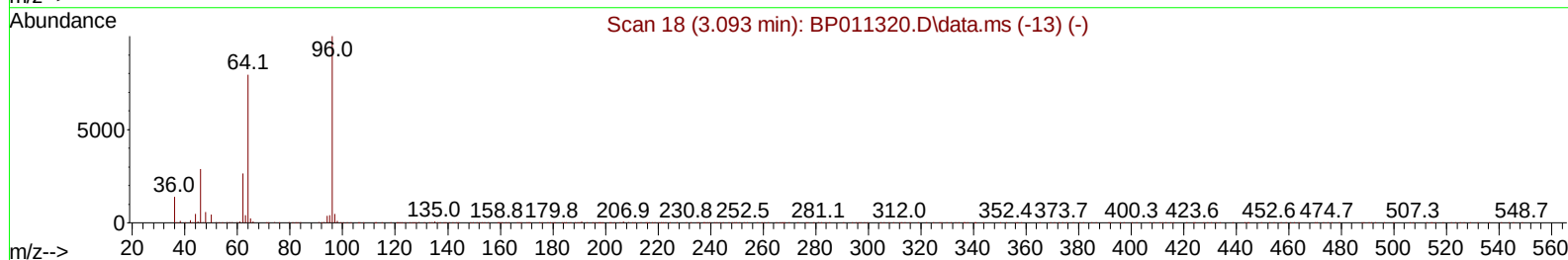
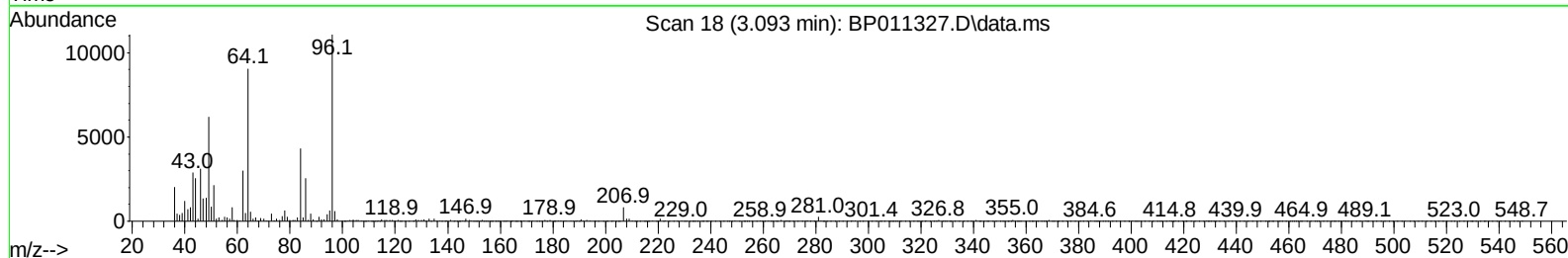
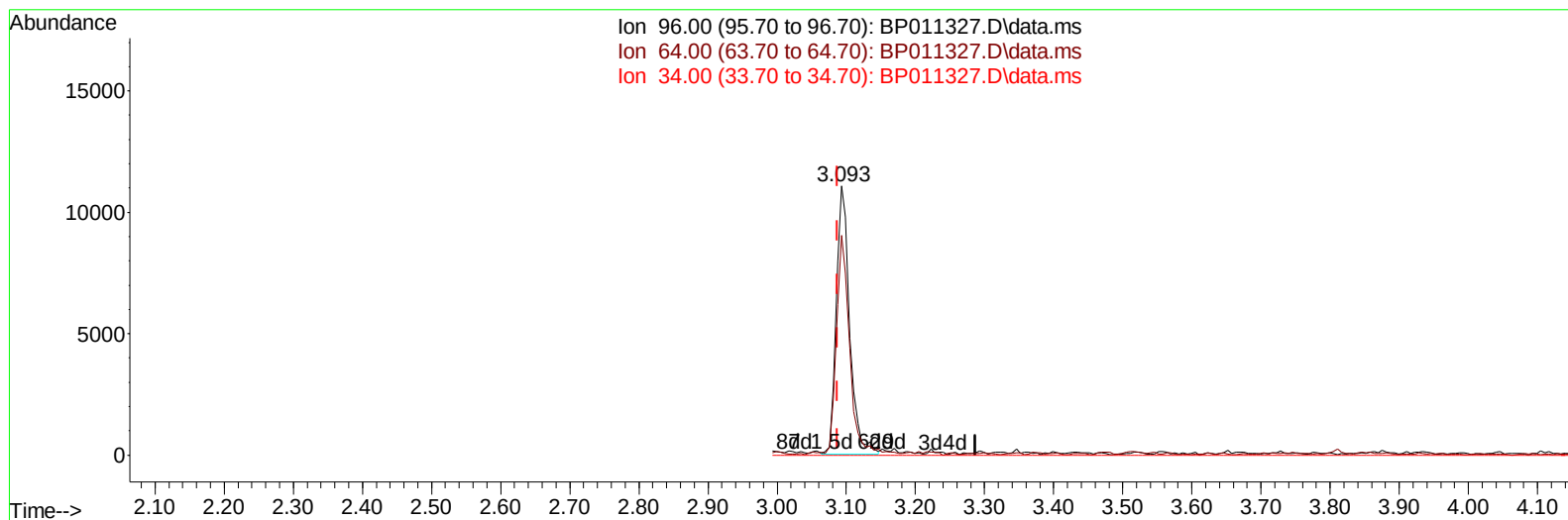
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TIC: BP011327.D\data.ms

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

3.093min (+ 0.006) 3.63 ng/uL m

response 14923

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	81.79#
34.00	0.00	0.00
0.00	0.00	0.00

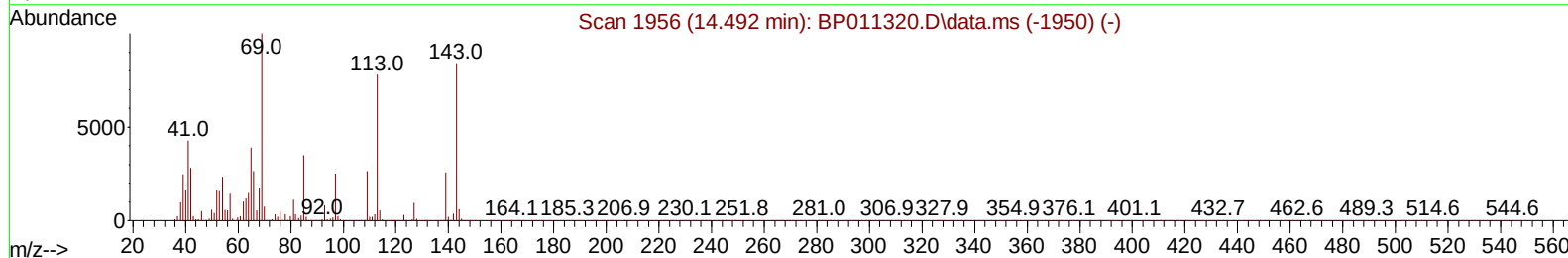
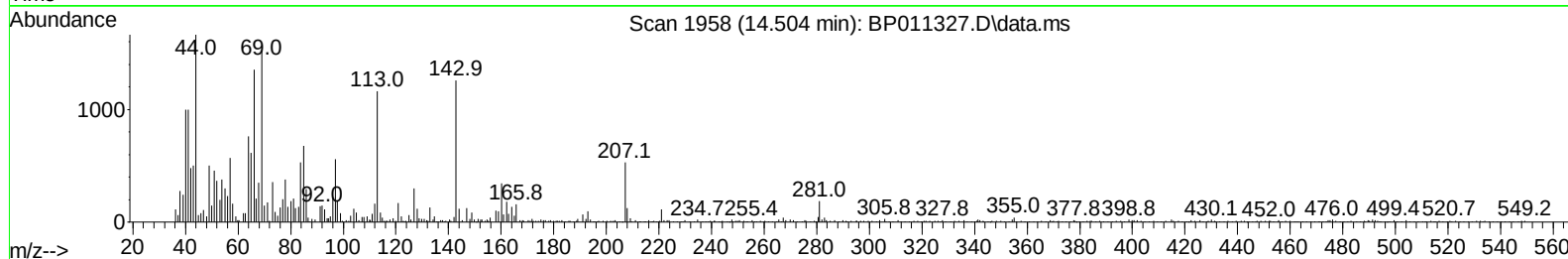
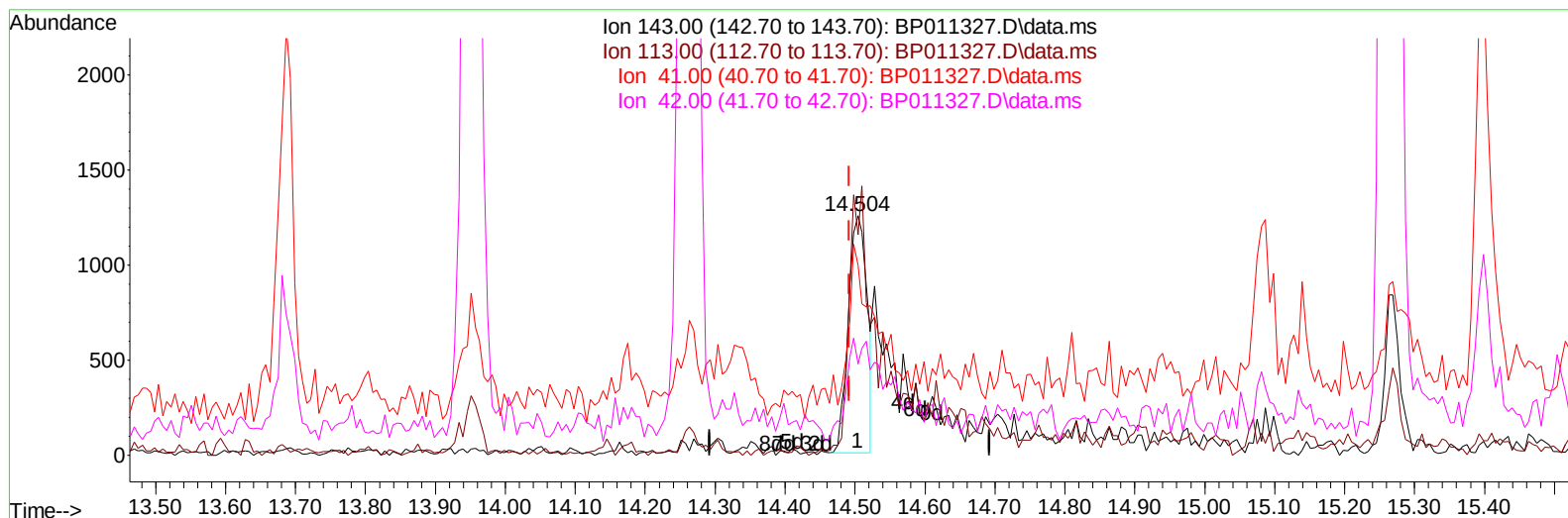
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TIC: BP011327.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.012) 0.63 ng/u1

response 2356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	91.98#
41.00	54.40	79.21#
42.00	35.50	38.10

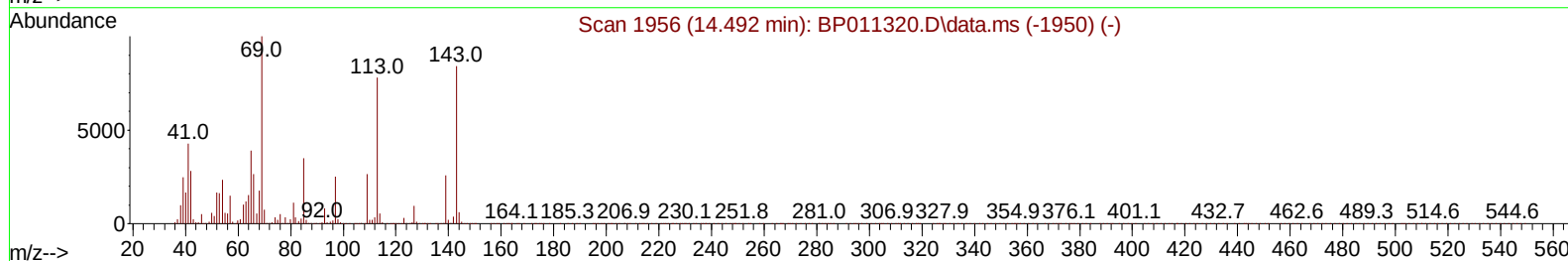
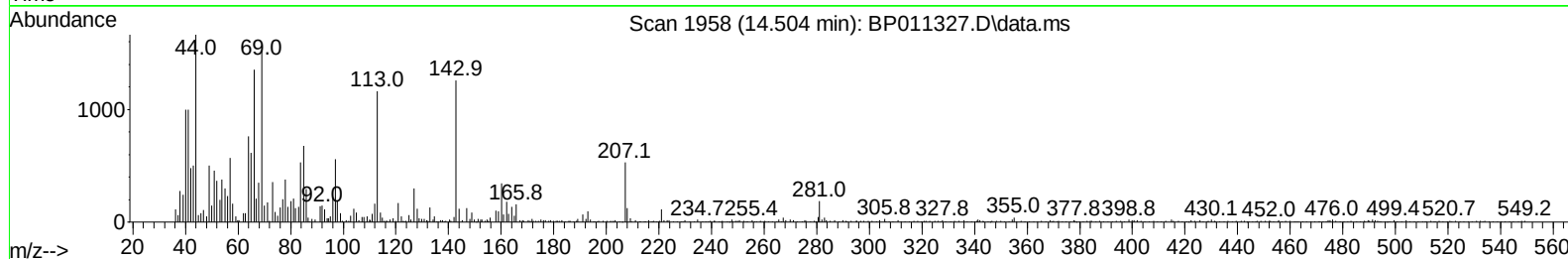
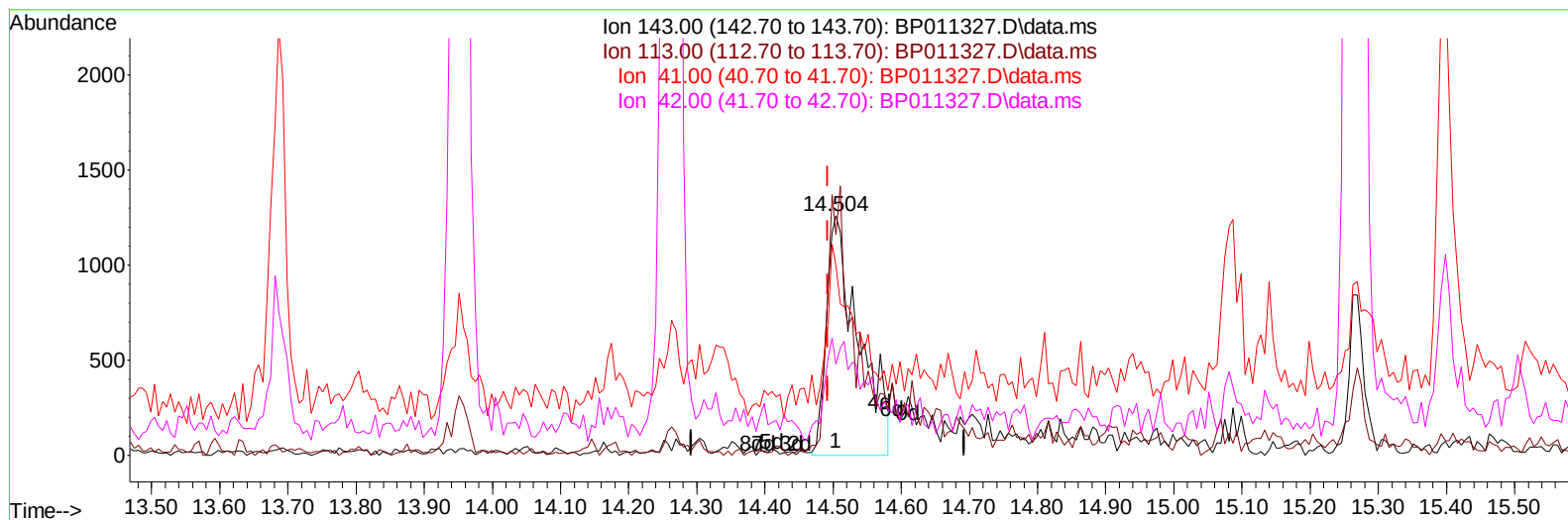
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Instrument :
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TIC: BP011327.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.012) 1.10 ng/ul m

response 4128

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	91.98#
41.00	54.40	79.21#
42.00	35.50	38.10

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	125946	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	521833	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	296336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	610483	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	576569	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	584005	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	14923m	3.632	ng/uL	0.00
4) Pyridine-d5	3.511	84	60653	5.853	ng/ul	0.01
7) Phenol-d5	6.775	99	31118	2.792	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	171583	21.711	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	88579	10.129	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	69132	7.899	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	87711	24.697	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	40562	12.475	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	86507	13.174	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	137559	11.983	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	610818	29.707	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	682327	27.568	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	4128m	1.099	ng/ul	0.01
60) Fluorene-d10	15.269	176	516426	29.369	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	25870	9.123	ng/ul	0.00
73) Anthracene-d10	17.134	188	874019	31.852	ng/ul	0.00
81) Pyrene-d10	19.398	212	1003707	31.598	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	917099	31.139	ng/ul	0.00

Target Compounds Qvalue

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

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