

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
Data File : BP021800.D
Acq On : 03 Sep 2024 23:30
Operator : RC/JU
Sample : P3734-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

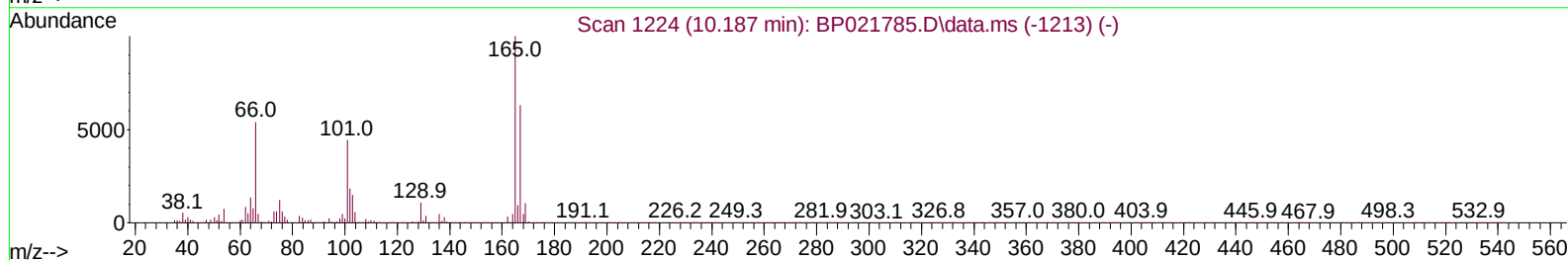
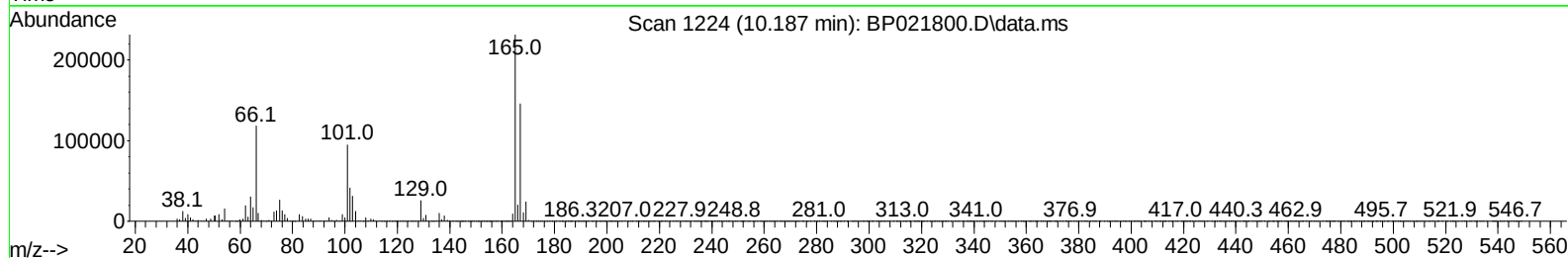
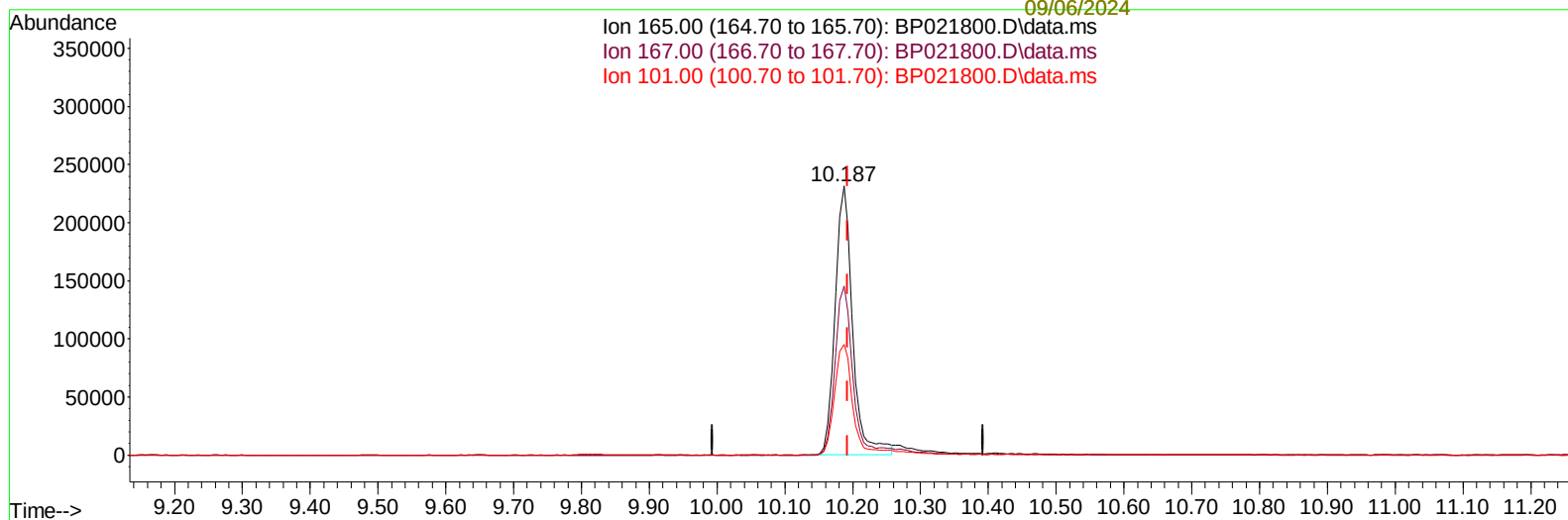
Instrument :
BNA_P
ClientSampleId :
A44B0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/04/2024
Supervised By :mohammad ahmed 09/07/2024

Supervised By :mohammad
ahmed
09/06/2024

Quant Time: Sep 04 04:04:12 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 29 23:56:59 2024
Response via : Initial Calibration



TIC: BP021800.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.187min (-0.006) 23.59 ng/ul

response 416111

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.93
101.00	41.40	41.19
0.00	0.00	0.00

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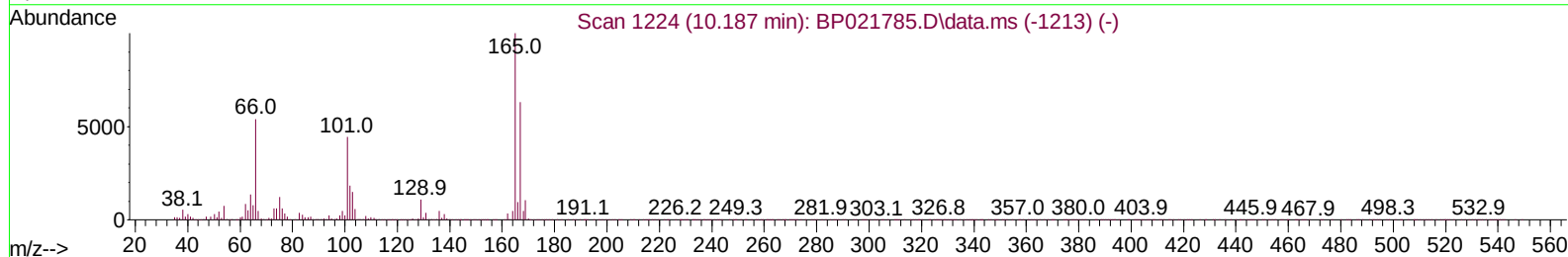
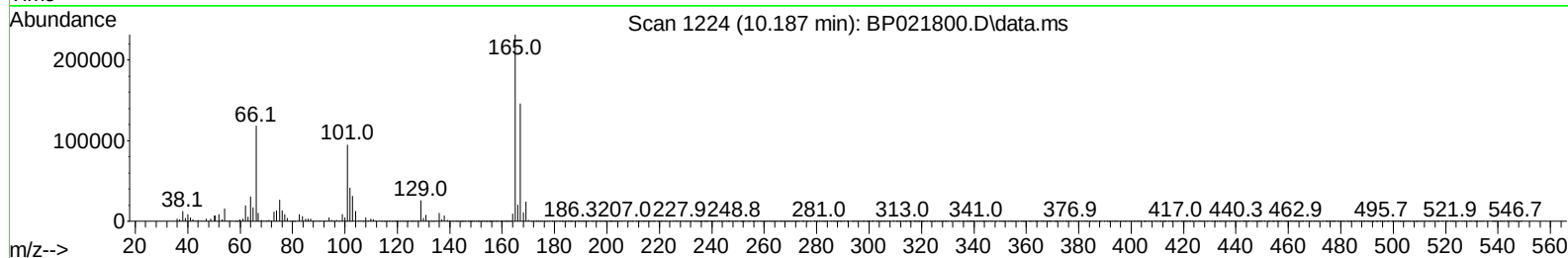
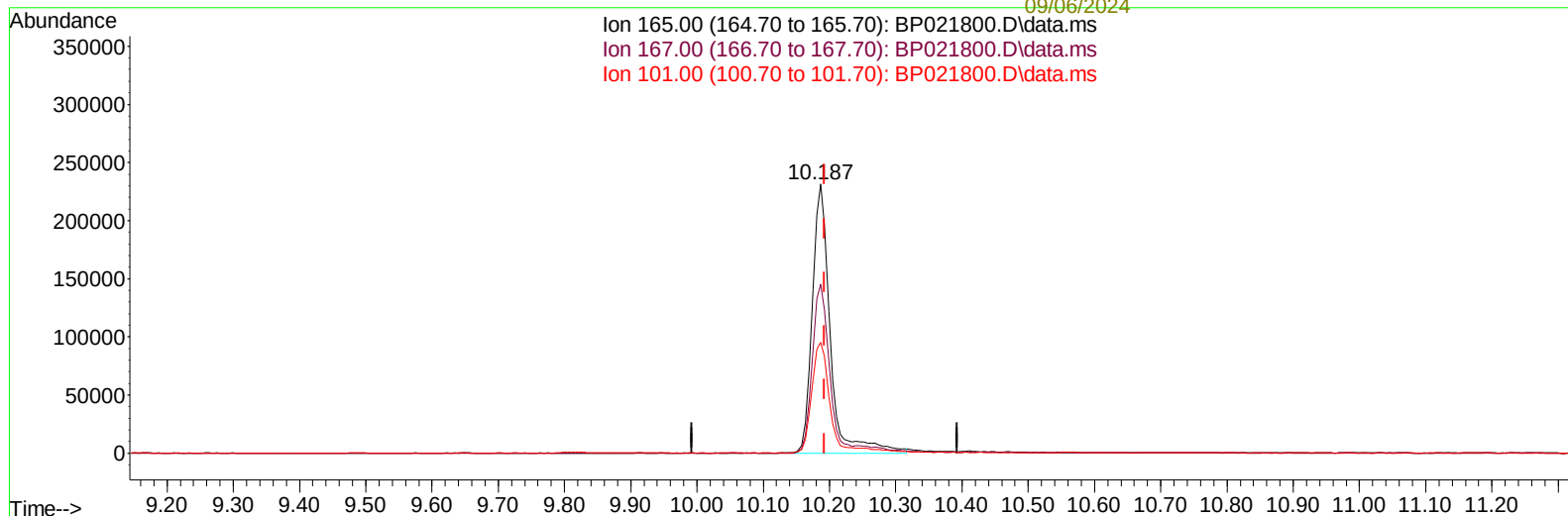
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TIC: BP021800.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.187min (-0.006) 24.70 ng/ul m

response 435702

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.93
101.00	41.40	41.19
0.00	0.00	0.00

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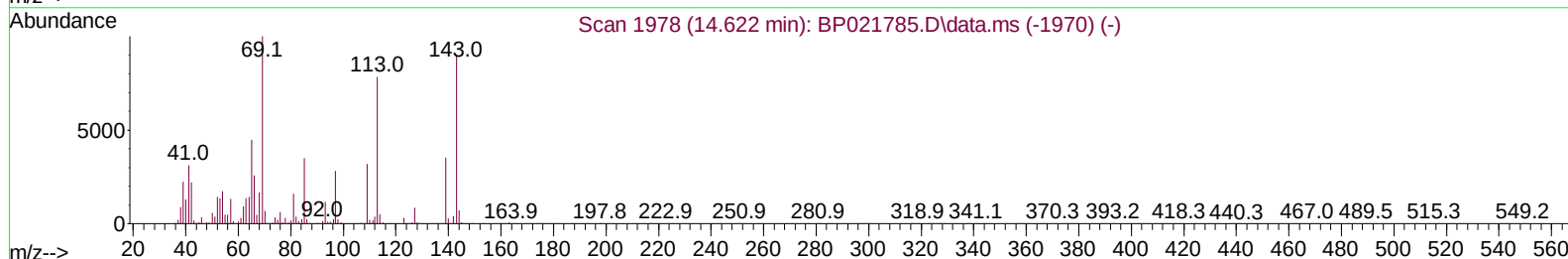
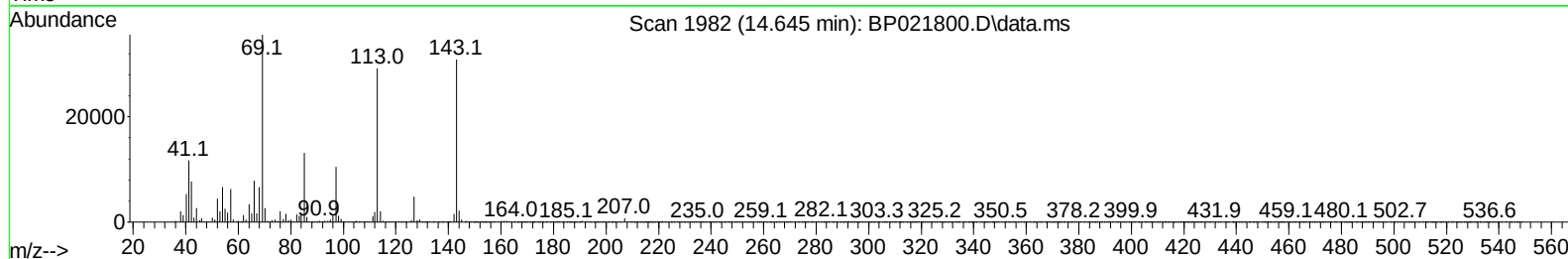
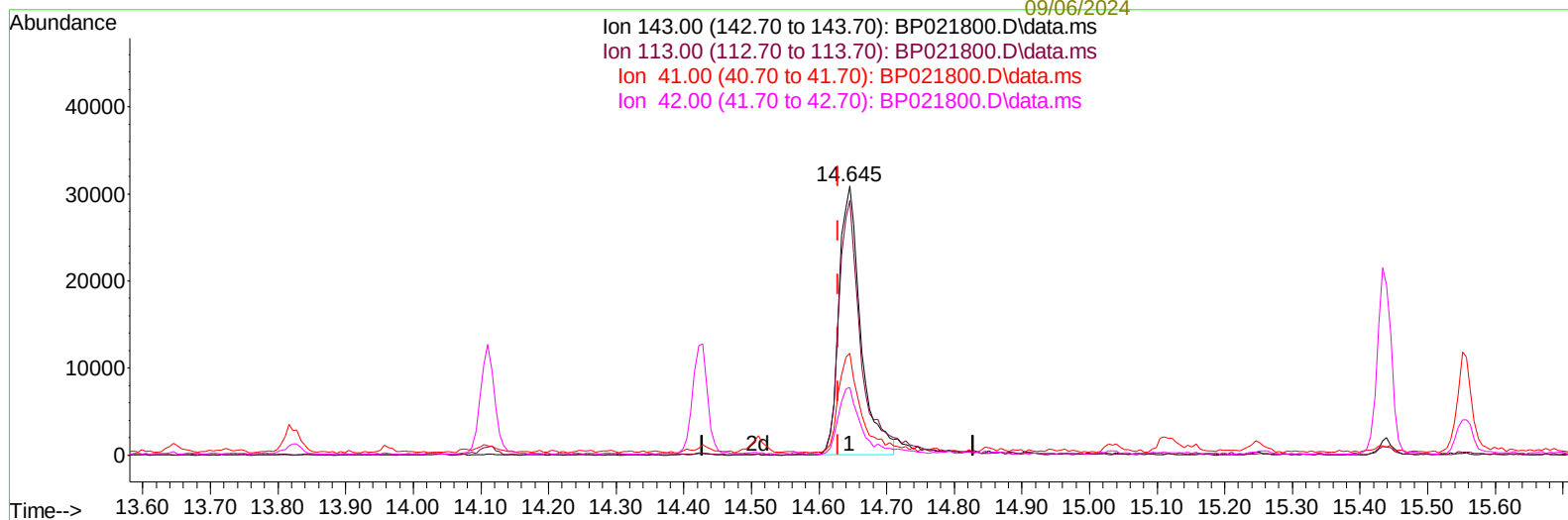
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TIC: BP021800.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.018) 6.72 ng/ul

response 68840

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.10	94.59
41.00	33.90	37.76
42.00	23.00	25.04

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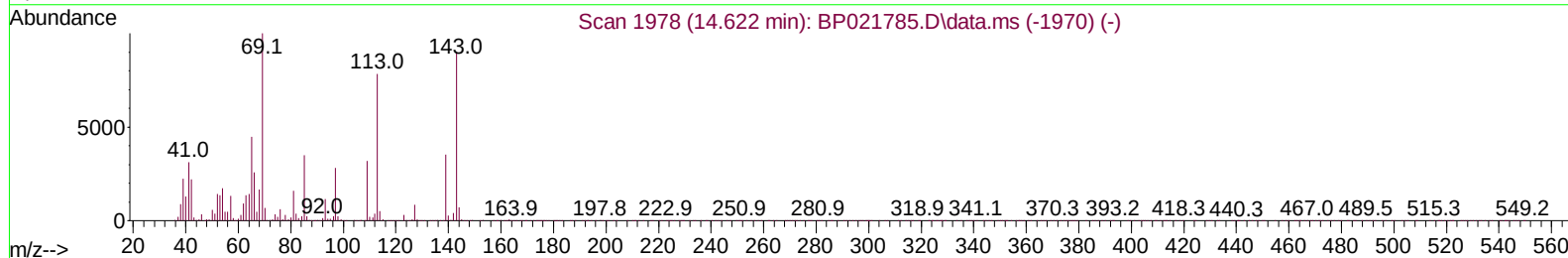
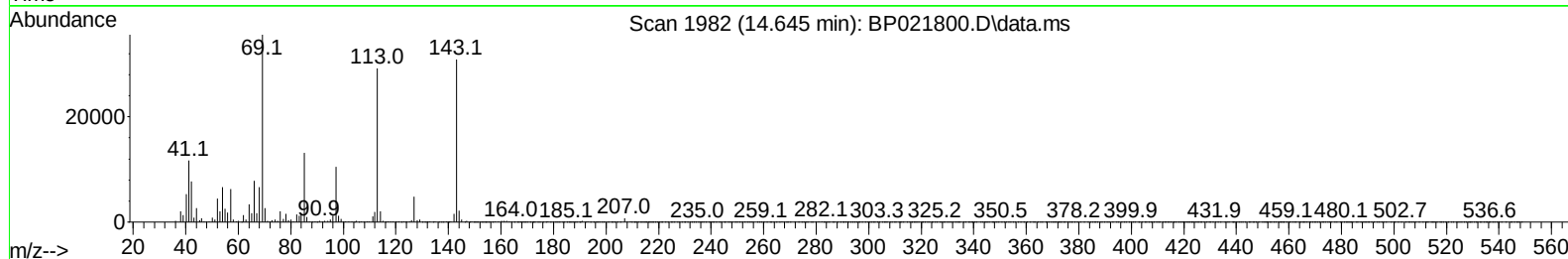
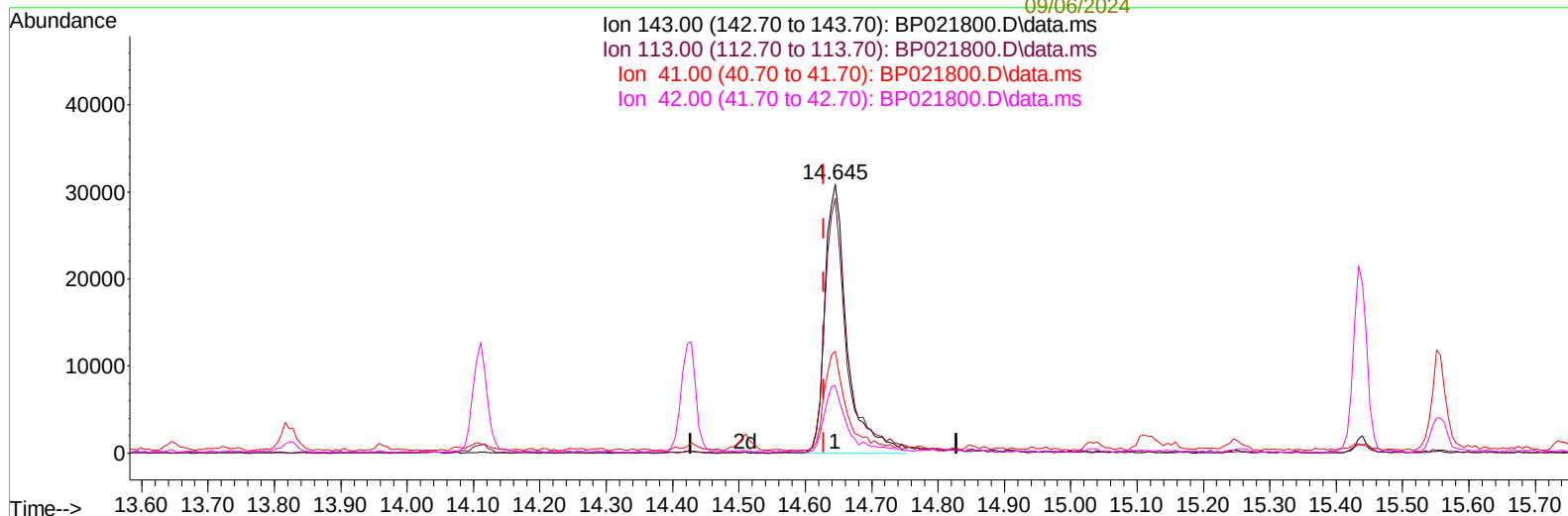
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TIC: BP021800.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.018) 6.99 ng/ul m

response 71644

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.10	94.59
41.00	33.90	37.76
42.00	23.00	25.04

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.775	152		276542	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.563	136		1096020	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164		704194	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188		1550850	20.000	ng/ul	0.00
79) Chrysene-d12	21.663	240		1637937	20.000	ng/ul	0.00
88) Perylene-d12	25.074	264		1867836	20.000	ng/ul	0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.275	96		32284	3.876	ng/uL	0.00
4) Pyridine-d5	3.687	84		150469	6.245	ng/ul	0.00
7) Phenol-d5	6.952	99		220823	7.482	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67		403462	25.121	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.316	132		447433	23.483	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113		370638	16.252	ng/ul	0.00
21) Nitrobenzene-d5	8.922	128		235833	26.745	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.646	143		241108	26.937	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165		435702m	24.702	ng/ul	0.00
31) 4-Chloroaniline-d4	10.693	131		145690	5.677	ng/ul	0.00
46) Dimethylphthalate-d6	13.822	166		1529477	28.385	ng/ul	-0.02
49) Acenaphthylene-d8	14.110	160		1633474	27.001	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.645	143		71644m	6.989	ng/ul	0.02
60) Fluorene-d10	15.440	176		1326947	29.292	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200		199487	23.387	ng/ul	0.00
73) Anthracene-d10	17.334	188		2281623	31.000	ng/ul	0.00
81) Pyrene-d10	19.692	212		2857401	30.518	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.827	264		3202718	32.156	ng/ul	-0.01
Target Compounds							
59) Diethylphthalate	15.263	149		56139	1.036	ng/ul	Qvalue 97

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

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