

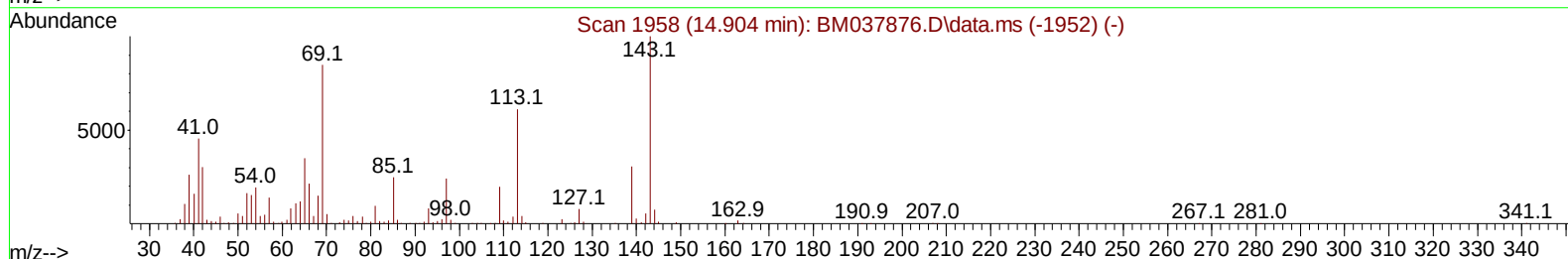
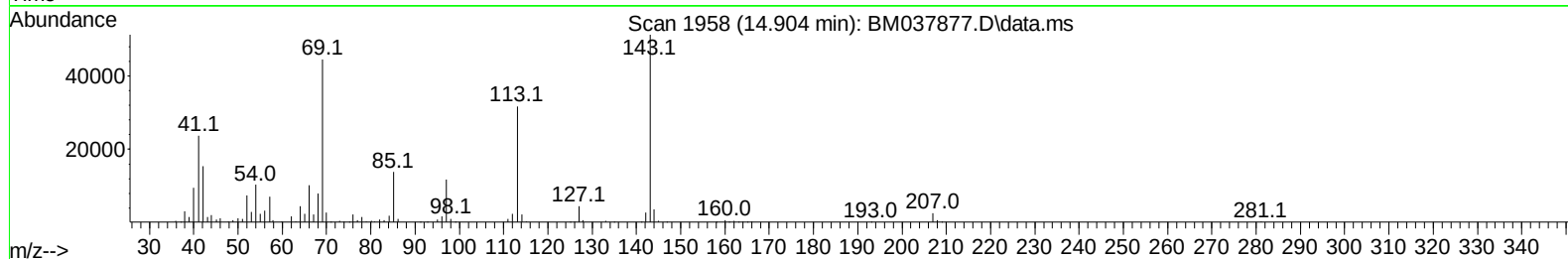
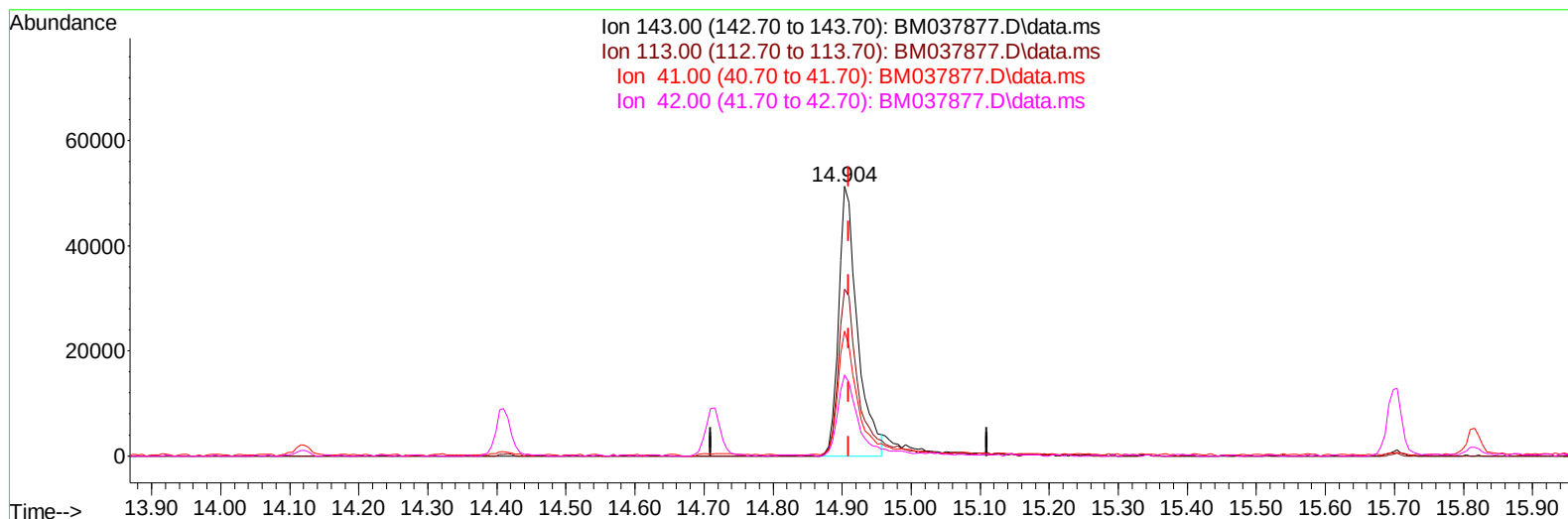
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120822\
Data File : BM037877.D
Acq On : 08 Dec 2022 09:08
Operator : CG/JU
Sample : PB149447BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK447

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/09/2022
Supervised By :mohammad ahmed 12/10/2022

Quant Time: Dec 08 10:20:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 08 01:49:06 2022
Response via : Initial Calibration



TIC: BM037877.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.904min (-0.006) 18.40 ng/ul****response 96667**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.30	61.90
41.00	43.90	46.36
42.00	30.30	30.04

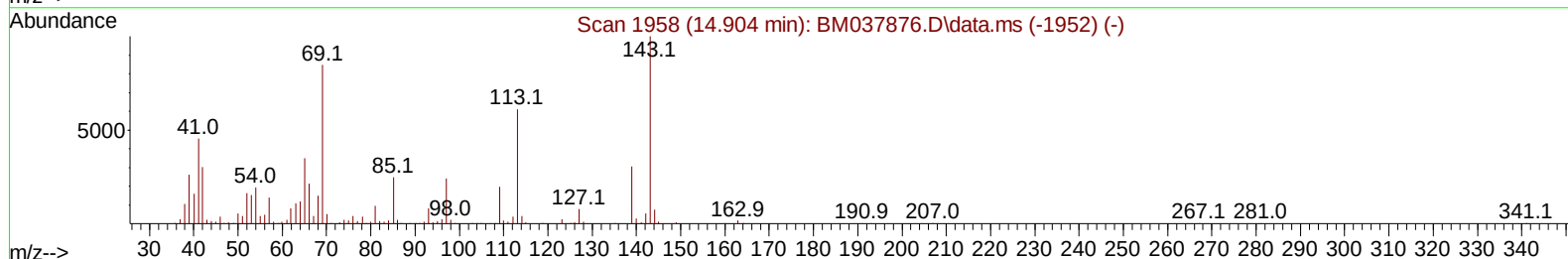
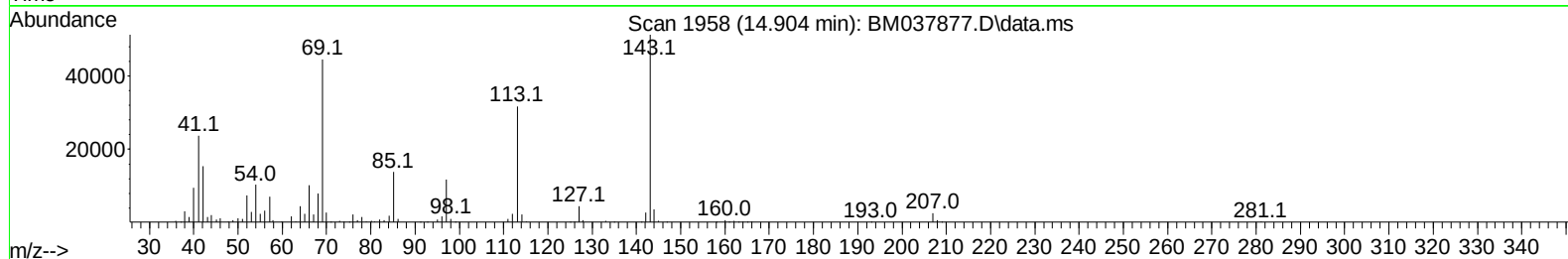
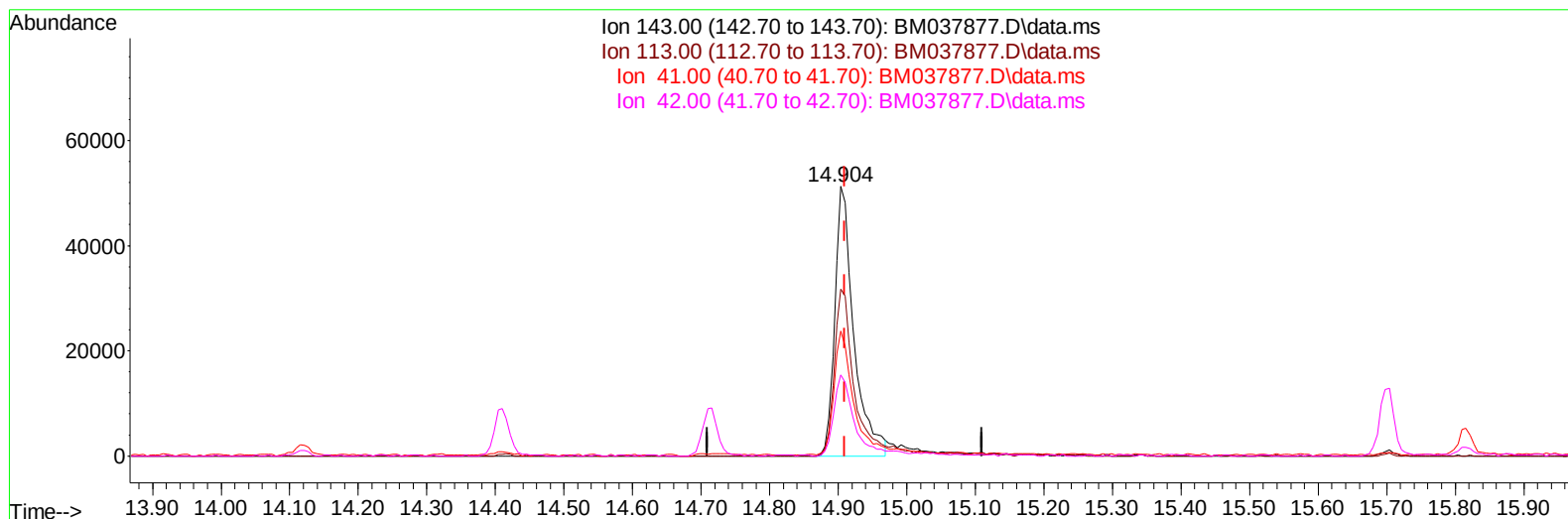
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TIC: BM037877.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (-0.006) 18.86 ng/ul m

response 99069

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.30	61.90
41.00	43.90	46.36
42.00	30.30	30.04

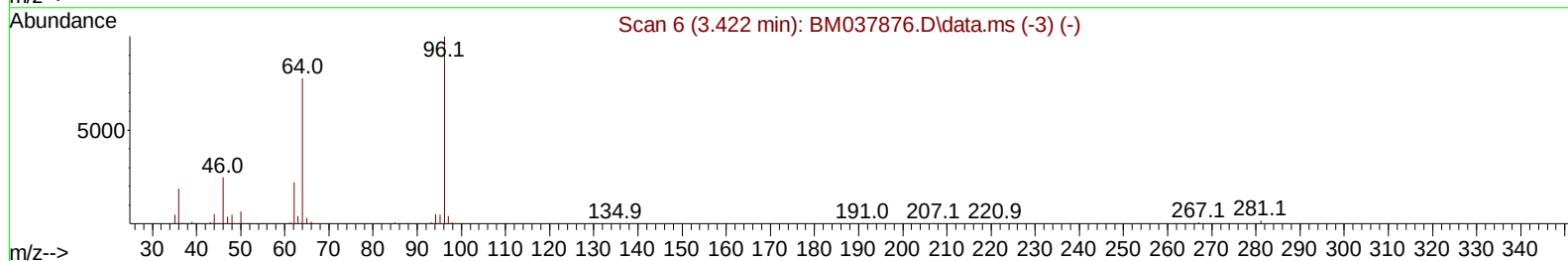
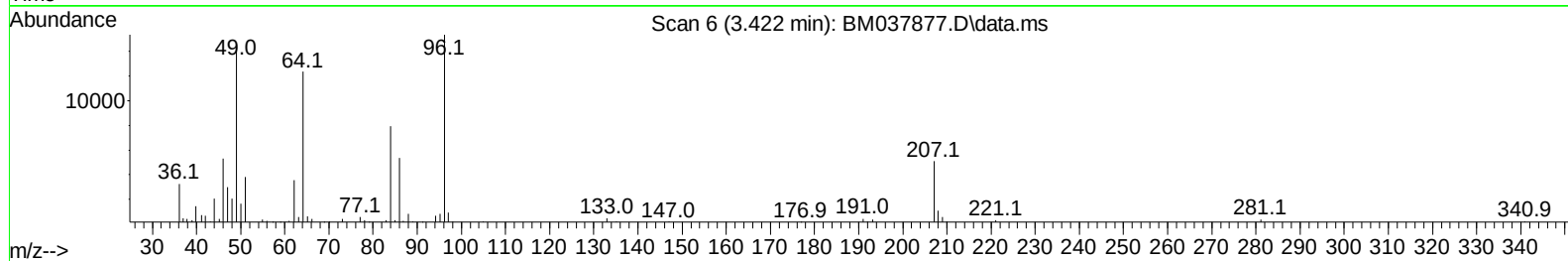
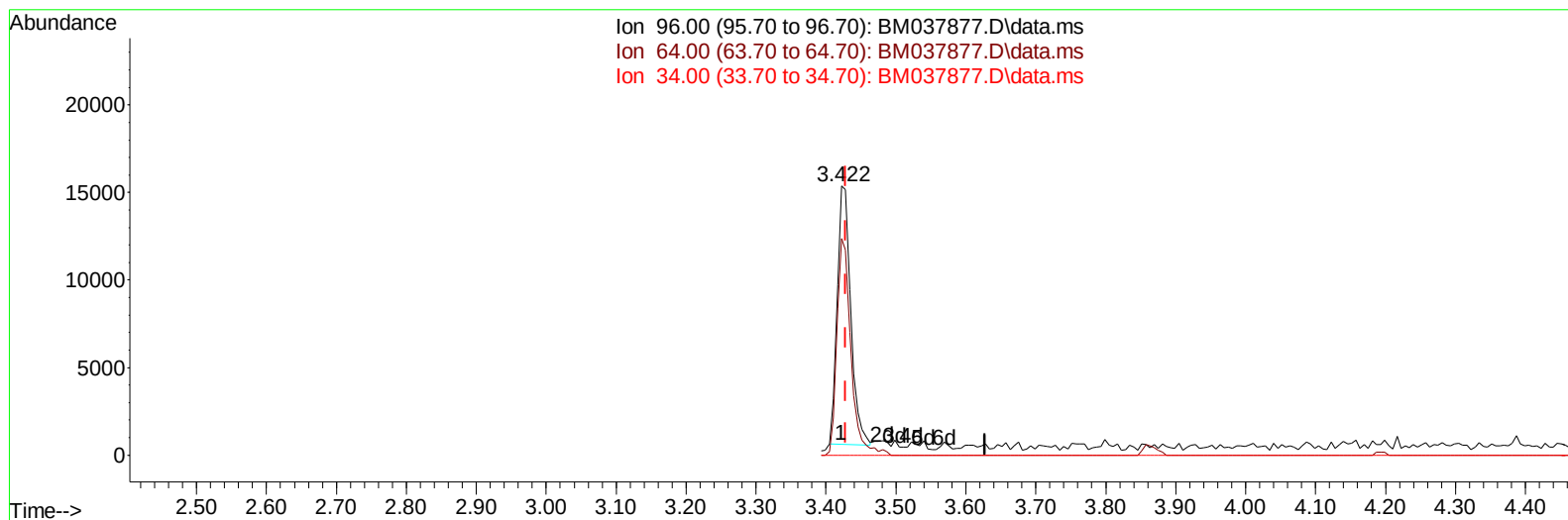
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 Misc :
 ALS Vial : 3 Sample Multiplier: 1

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TIC: BM037877.D\data.ms

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

3.422min (-0.006) 4.77 ng/uL

response 20176

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	80.41
34.00	0.00	0.00
0.00	0.00	0.00

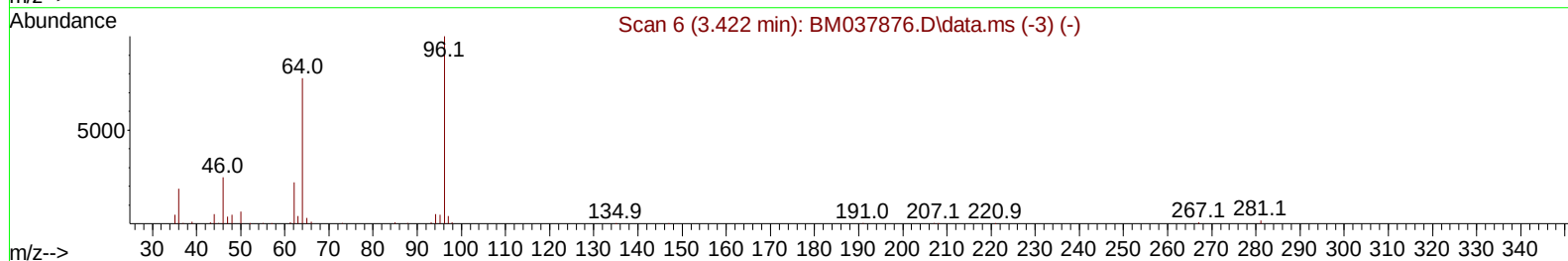
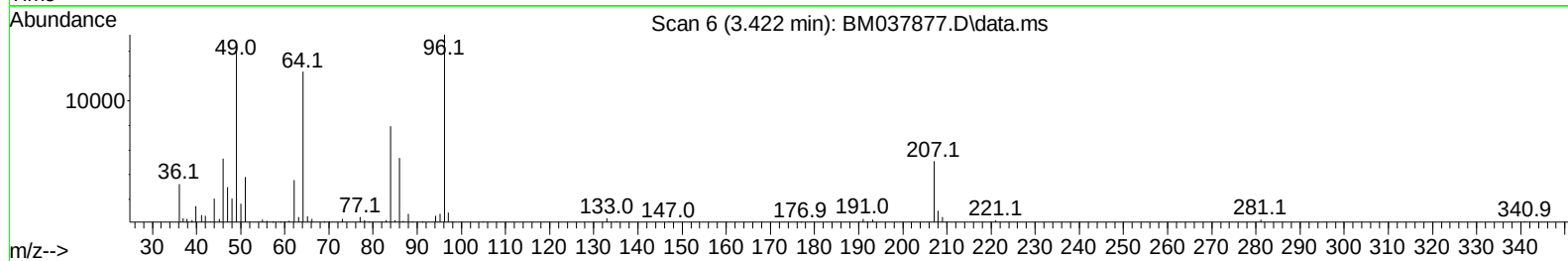
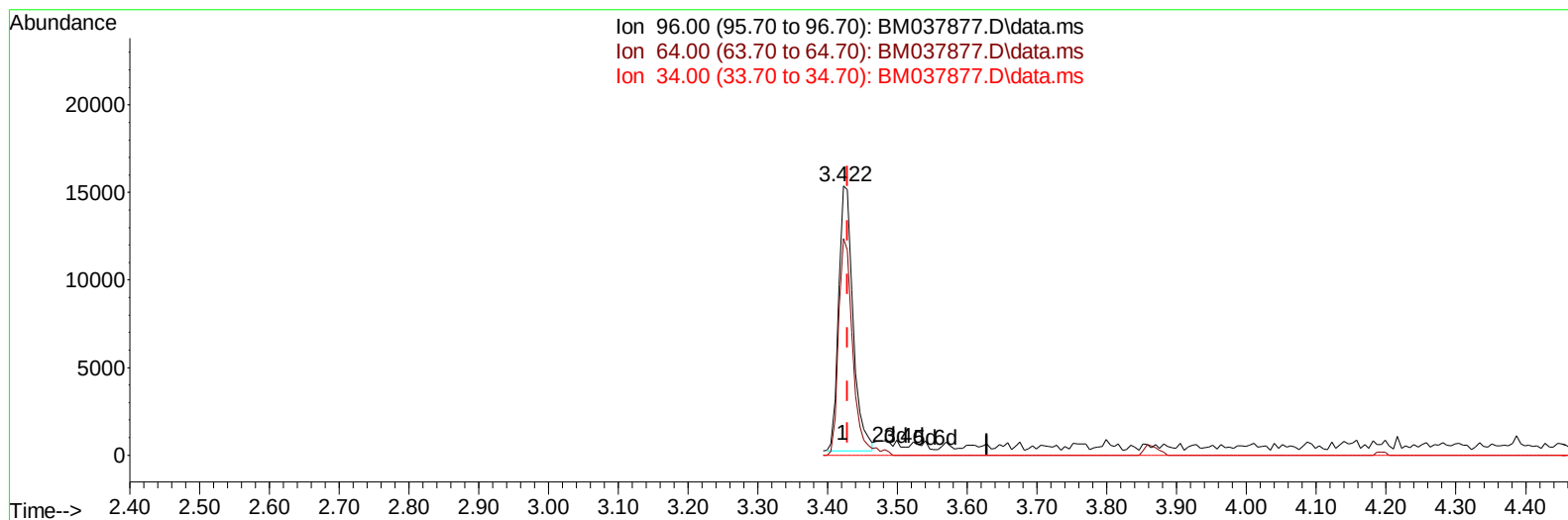
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 Operator : CG/JU
 Sample : PB149447BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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TIC: BM037877.D\data.ms

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

3.422min (-0.006) 5.11 ng/uL m

response 21612

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	80.41
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	192092	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	778215	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	432298	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	792138	20.000	ng/ul	0.00
79) Chrysene-d12	21.621	240	668582	20.000	ng/ul	0.00
88) Perylene-d12	24.103	264	724081	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	21612m	5.108	ng/uL	0.00
4) Pyridine-d5	3.863	84	323172	25.750	ng/ul	0.00
7) Phenol-d5	7.240	99	408183	26.150	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	254860	26.813	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	352691	27.543	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	321351	26.373	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128	171584	27.847	ng/ul	0.00
24) 2-Nitrophenol-d4	9.987	143	179447	27.791	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	313947	25.598	ng/ul	0.00
31) 4-Chloroaniline-d4	11.045	131	443874	26.219	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	848265	27.523	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	1017310	27.007	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	99069m	18.861	ng/ul	0.00
60) Fluorene-d10	15.704	176	745376	27.146	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	72210	17.050	ng/ul	0.00
73) Anthracene-d10	17.551	188	1067981	28.888	ng/ul	0.00
81) Pyrene-d10	19.833	212	1128811	28.107	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.944	264	1066210	29.481	ng/ul	0.00

Target Compounds Qvalue

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

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