

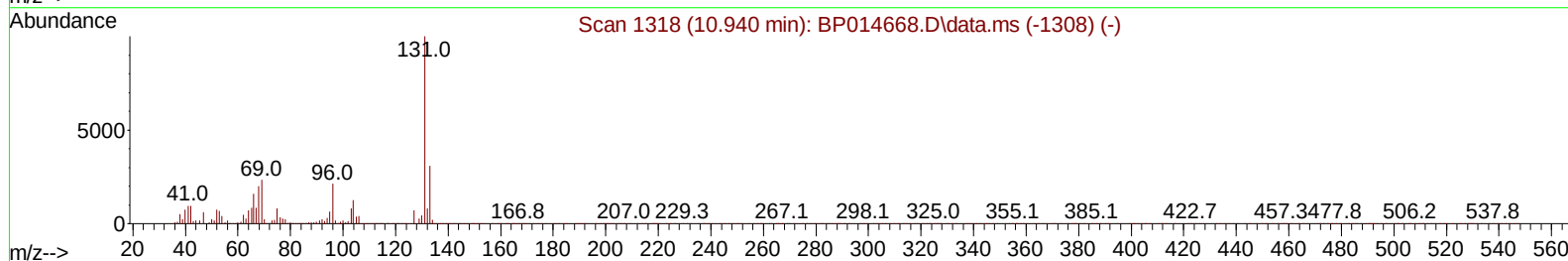
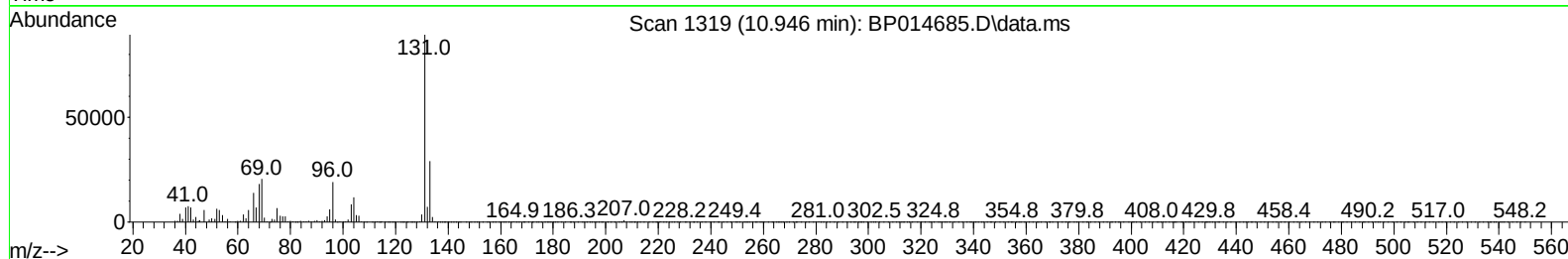
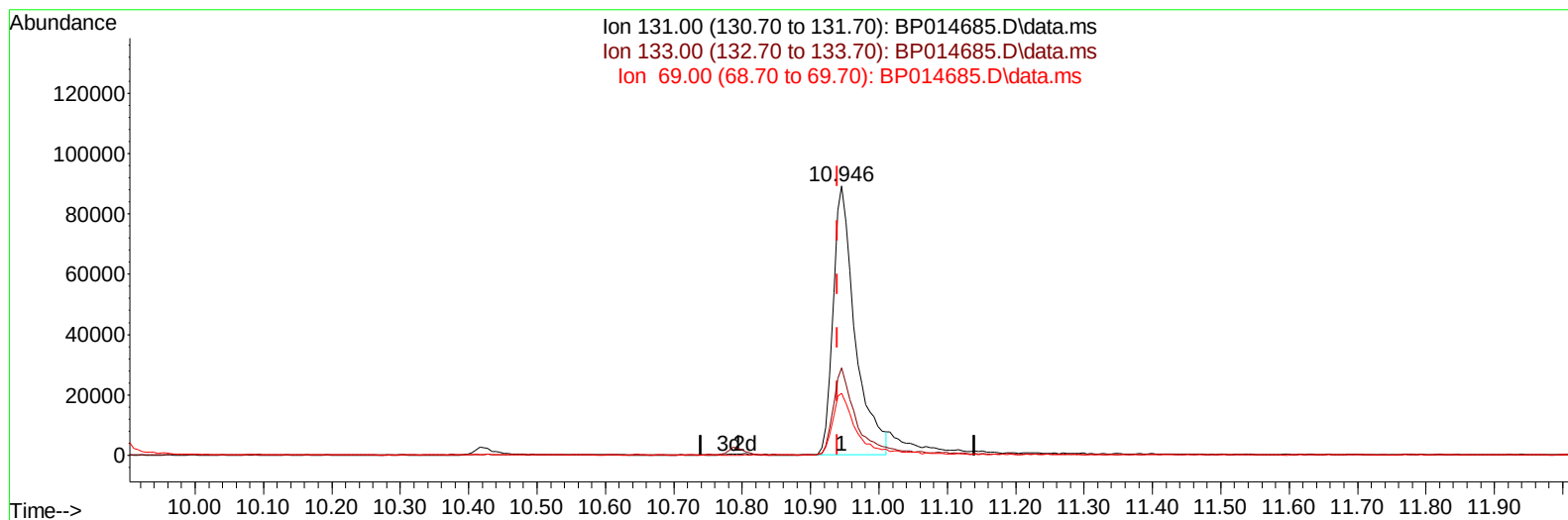
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041223\
 Data File : BP014685.D
 Acq On : 12 Apr 2023 22:24
 Operator : CG/JU
 Sample : PB152042BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK042

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:05:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 13 02:01:50 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/13/2023
 Supervised By :Jagrut Upadhyay 04/13/2023



TIC: BP014685.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.946min (+ 0.006) 19.21 ng/ul

response 199213

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.51
69.00	21.30	23.00
0.00	0.00	0.00

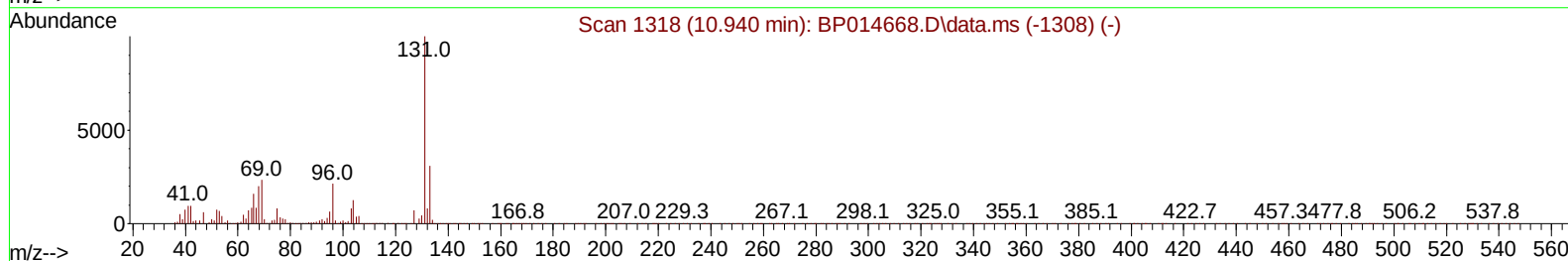
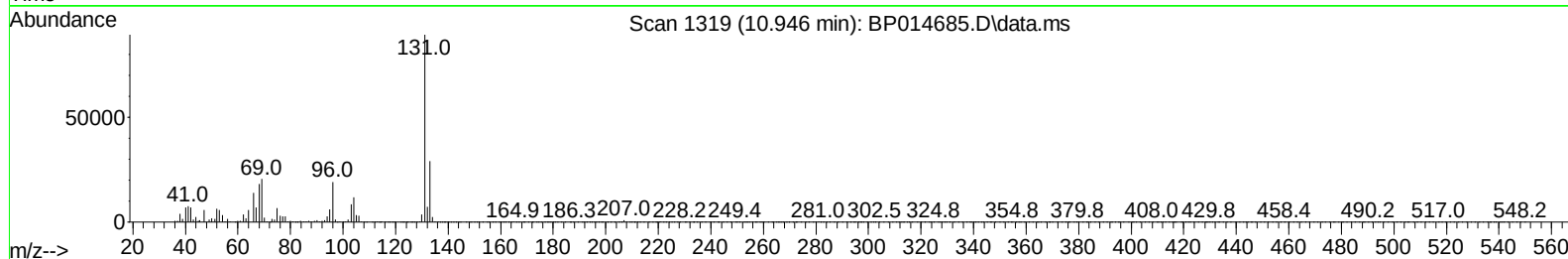
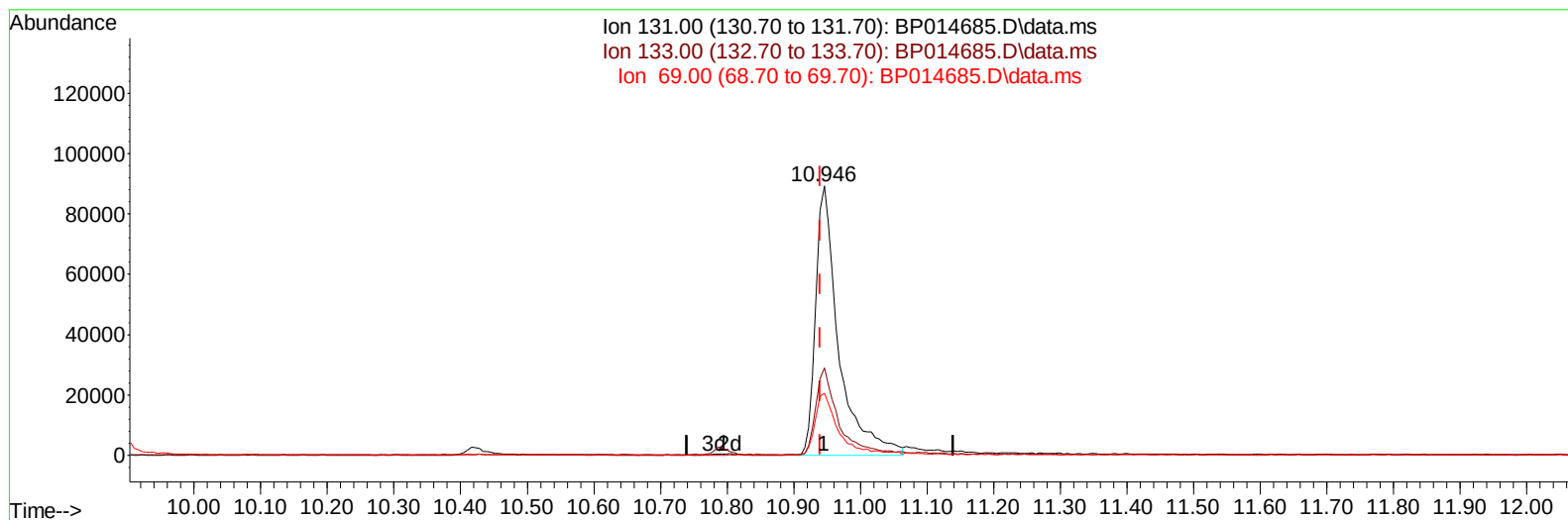
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TIC: BP014685.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.946min (+ 0.006) 20.62 ng/ul m

response 213864

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.51
69.00	21.30	23.00
0.00	0.00	0.00

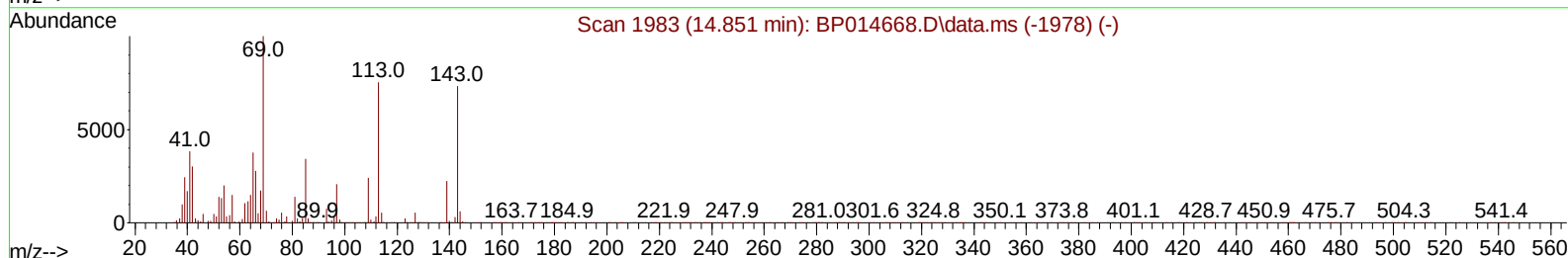
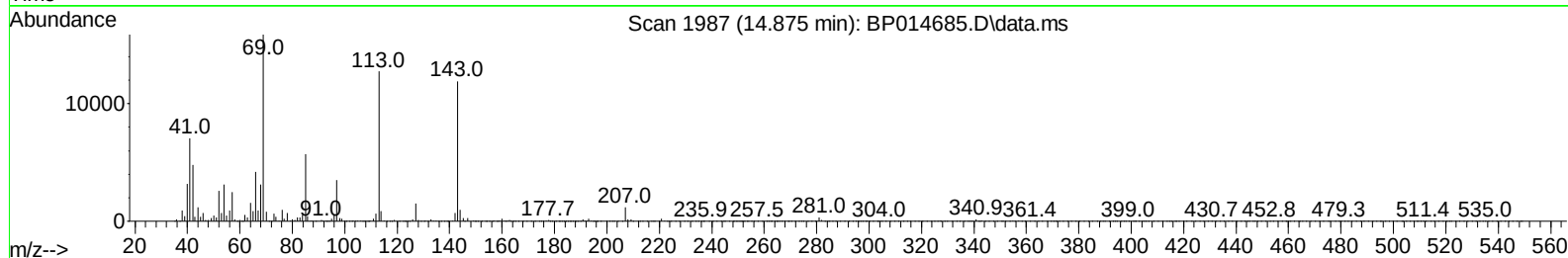
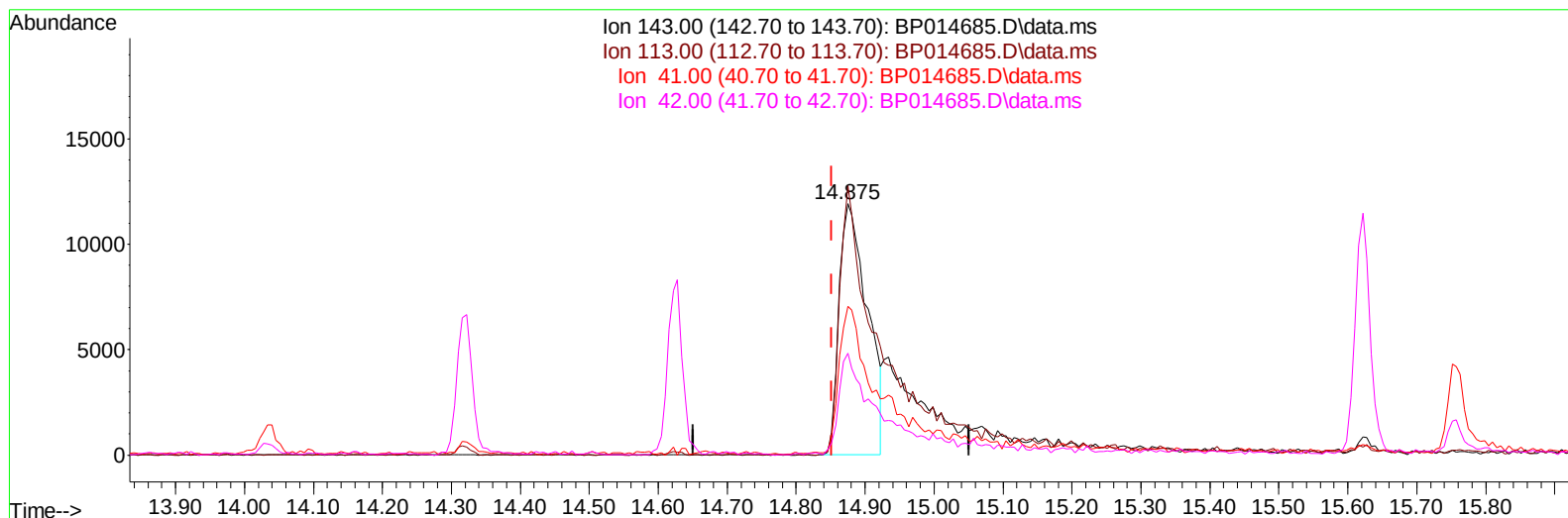
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014685.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.024) 10.83 ng/ul

response 33957

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	106.94
41.00	44.20	59.09#
42.00	31.50	40.38#

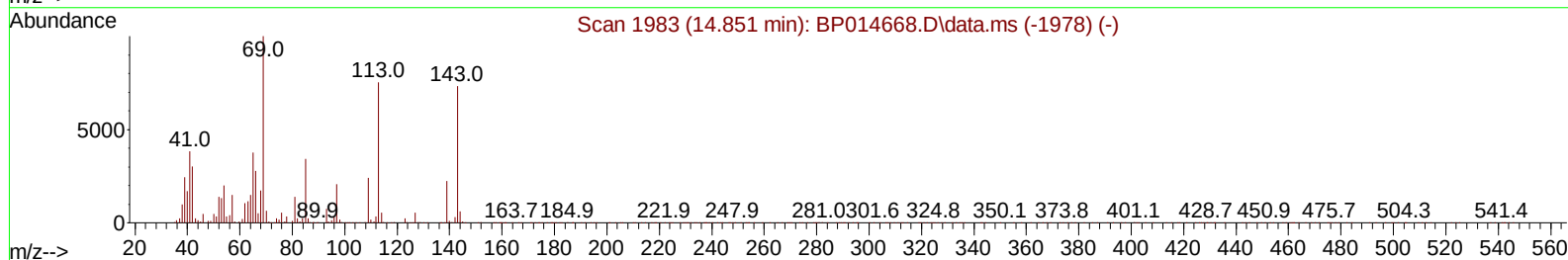
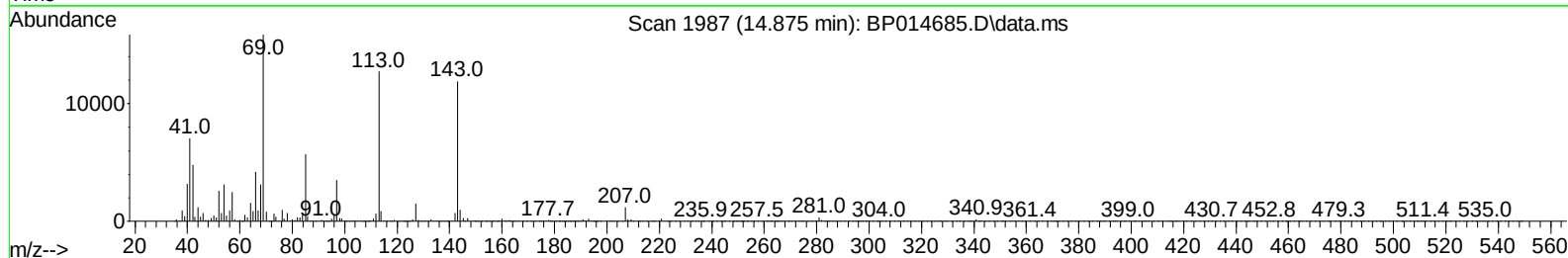
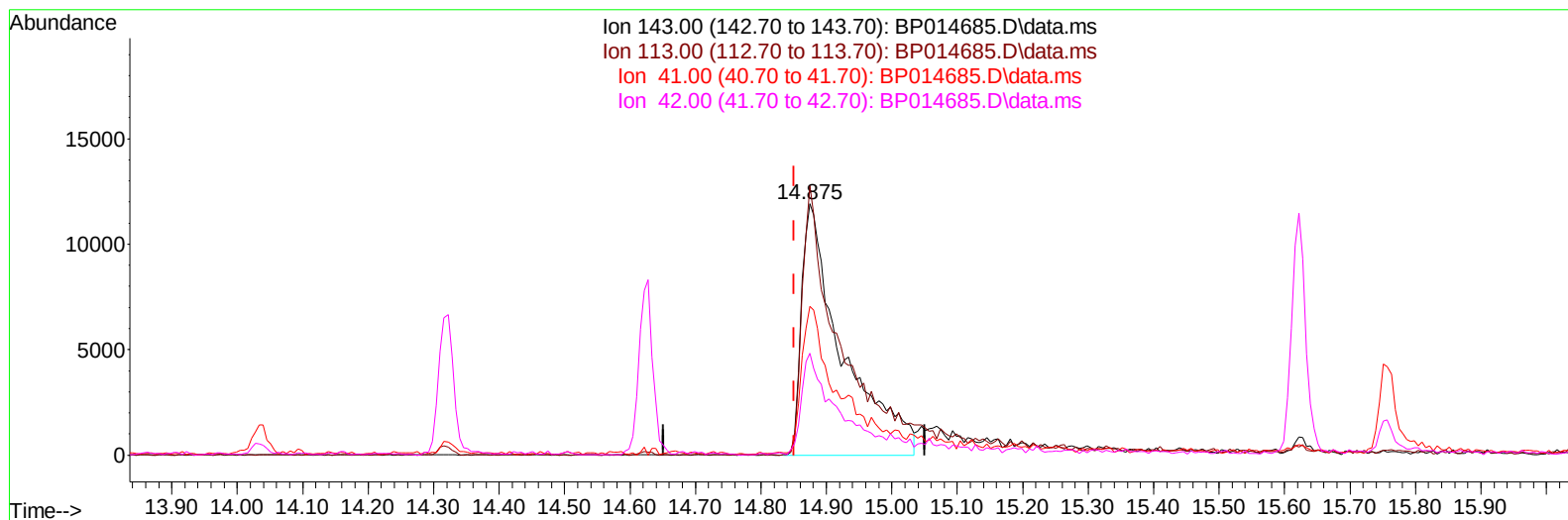
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Sample : PB152042BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP014685.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.024) 16.50 ng/ul m

response 51753

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	106.94
41.00	44.20	59.09#
42.00	31.50	40.38#

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 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	108749	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	467706	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.622	164	256084	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	530593	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	486909	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	581171	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	17066	6.086	ng/uL	0.00
4) Pyridine-d5	3.793	84	167926	23.367	ng/ul	0.00
7) Phenol-d5	7.140	99	252631	25.189	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	182994	28.438	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	196383	27.091	ng/ul	0.00
15) 4-Methylphenol-d8	8.693	113	197168	24.372	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	90762	24.038	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	105026	24.292	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	165562	21.027	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	213864m	20.620	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	520893	25.087	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	580592	25.200	ng/ul	0.00
54) 4-Nitrophenol-d4	14.875	143	51753m	16.498	ng/ul	0.02
60) Fluorene-d10	15.622	176	435013	25.123	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	71850	18.403	ng/ul	0.00
73) Anthracene-d10	17.486	188	659657	25.689	ng/ul	0.00
81) Pyrene-d10	19.722	212	755273	23.093	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	819412	26.538	ng/ul	0.00

Target Compounds Qvalue

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

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