

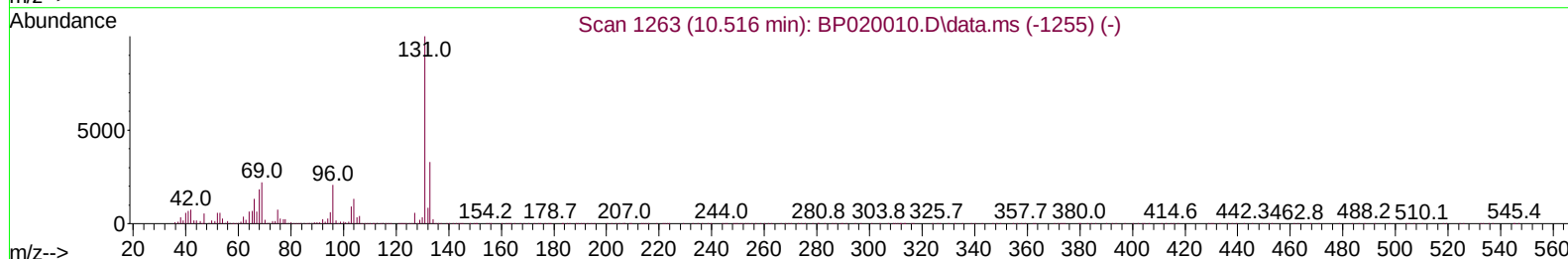
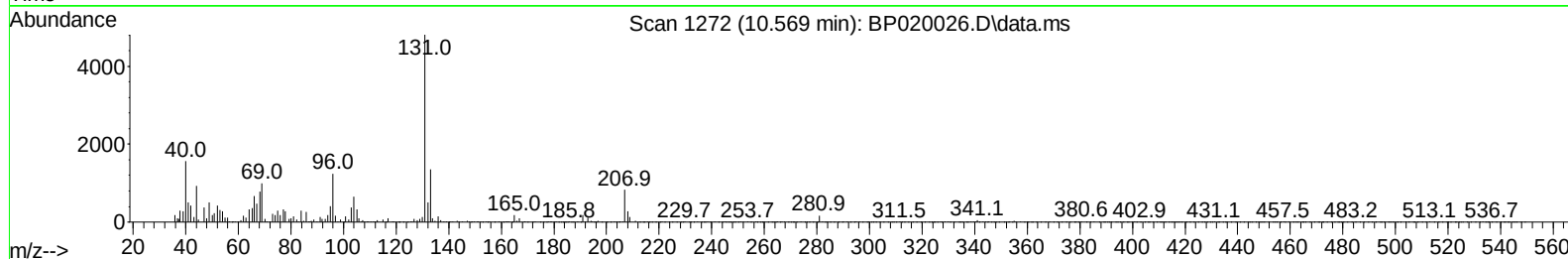
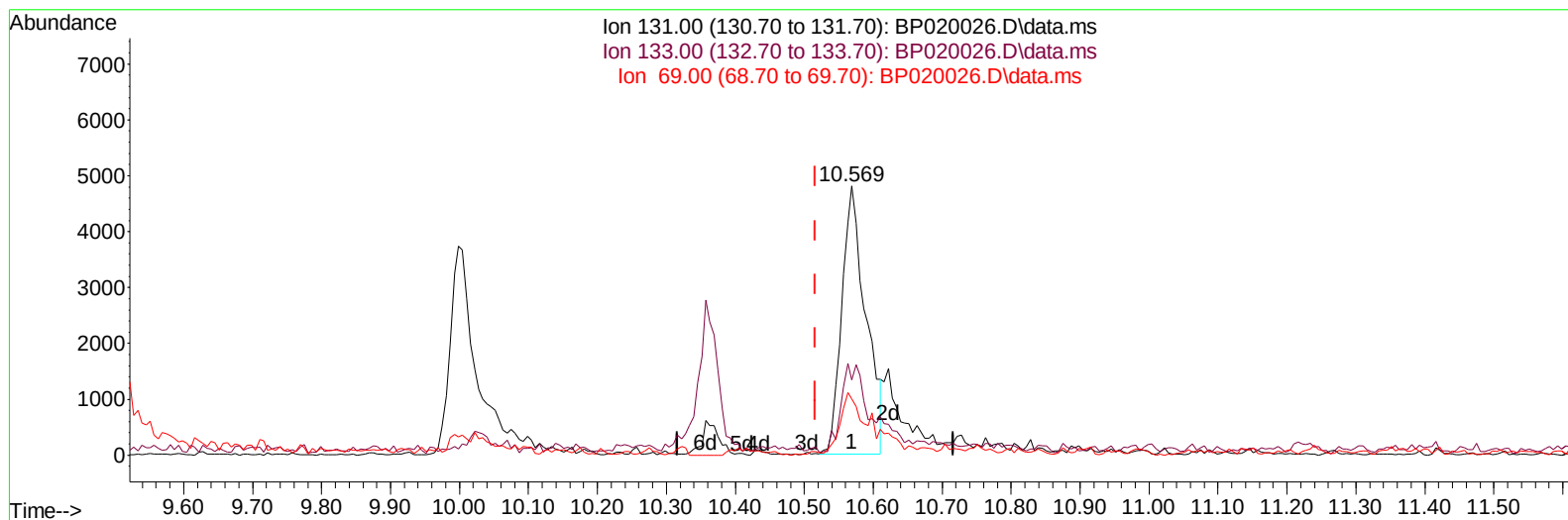
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
Data File : BP020026.D
Acq On : 29 Mar 2024 01:57
Operator : MA/JU
Sample : PB159808TB 10X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SLEB808

Manual IntegrationsAPPROVED

Quant Time: Mar 29 04:37:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 28 14:08:48 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
Supervised By :Jagrut Upadhyay 03/29/2024



TIC: BP020026.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.569min (+ 0.053) 0.96 ng/uI

response 11536

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	28.02
69.00	22.40	20.84
0.00	0.00	0.00

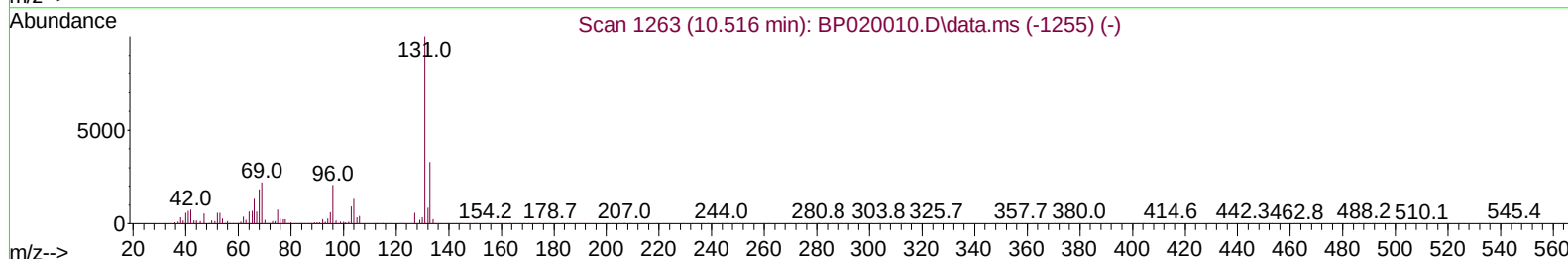
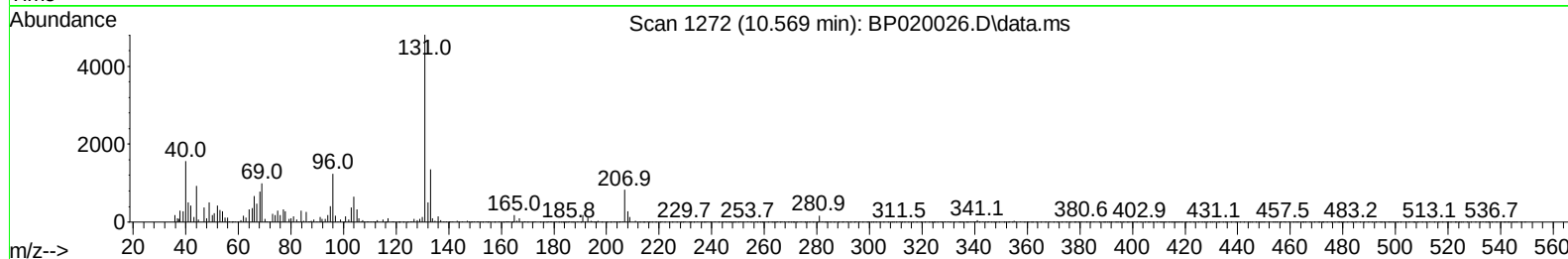
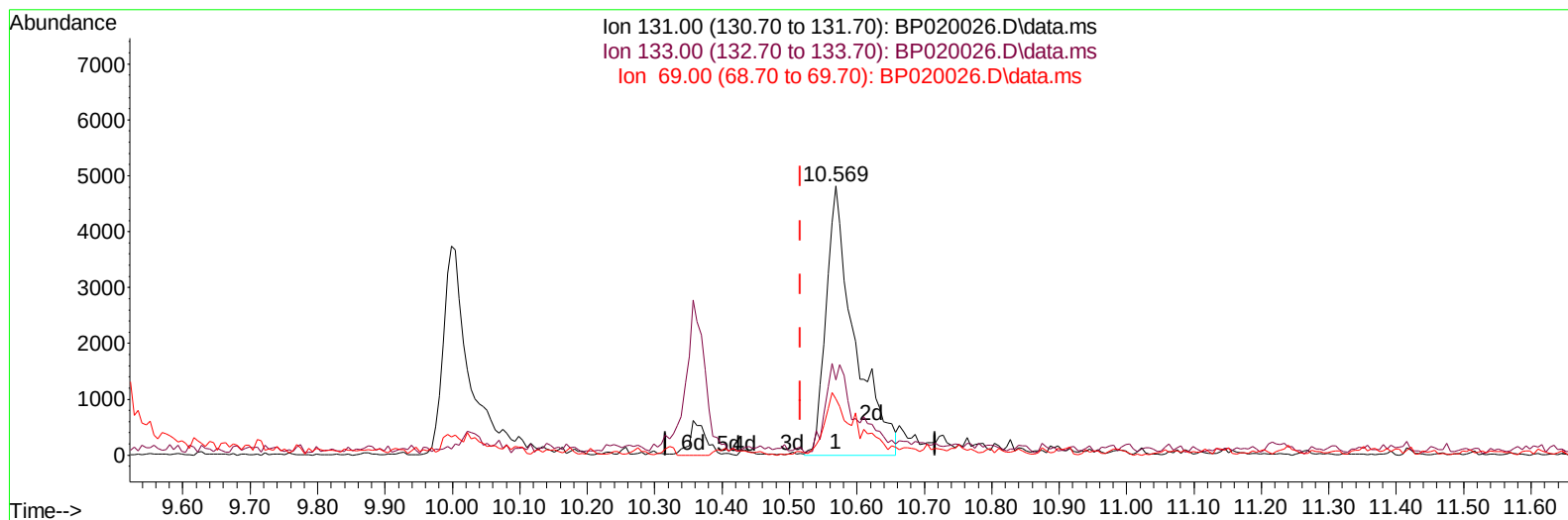
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(31) 4-Chloroaniline-d4 (S)

10.569min (+ 0.053) 1.17 ng/ul m

response 14080

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	28.02
69.00	22.40	20.84
0.00	0.00	0.00

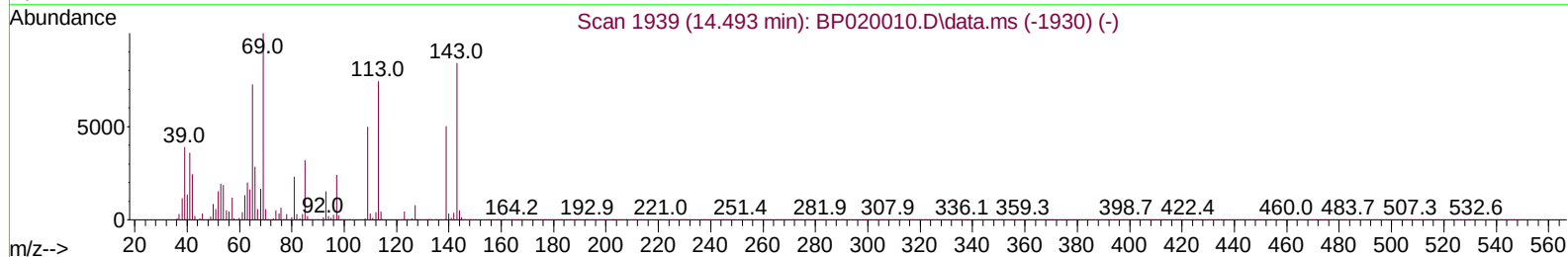
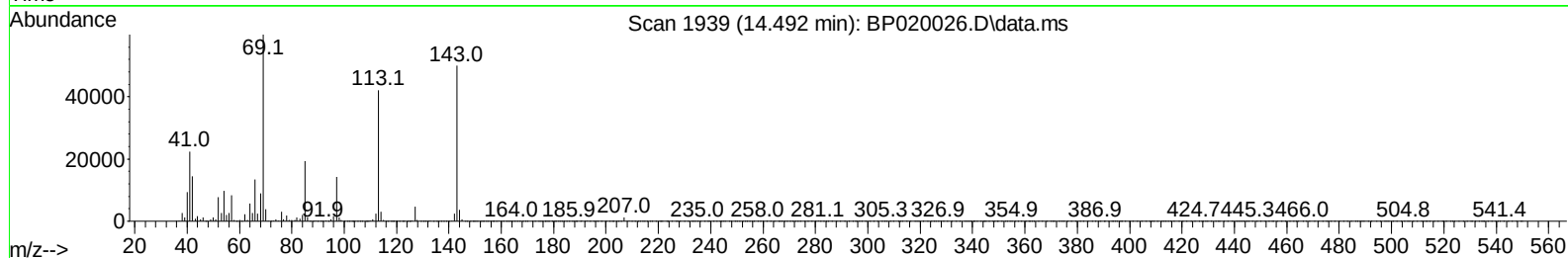
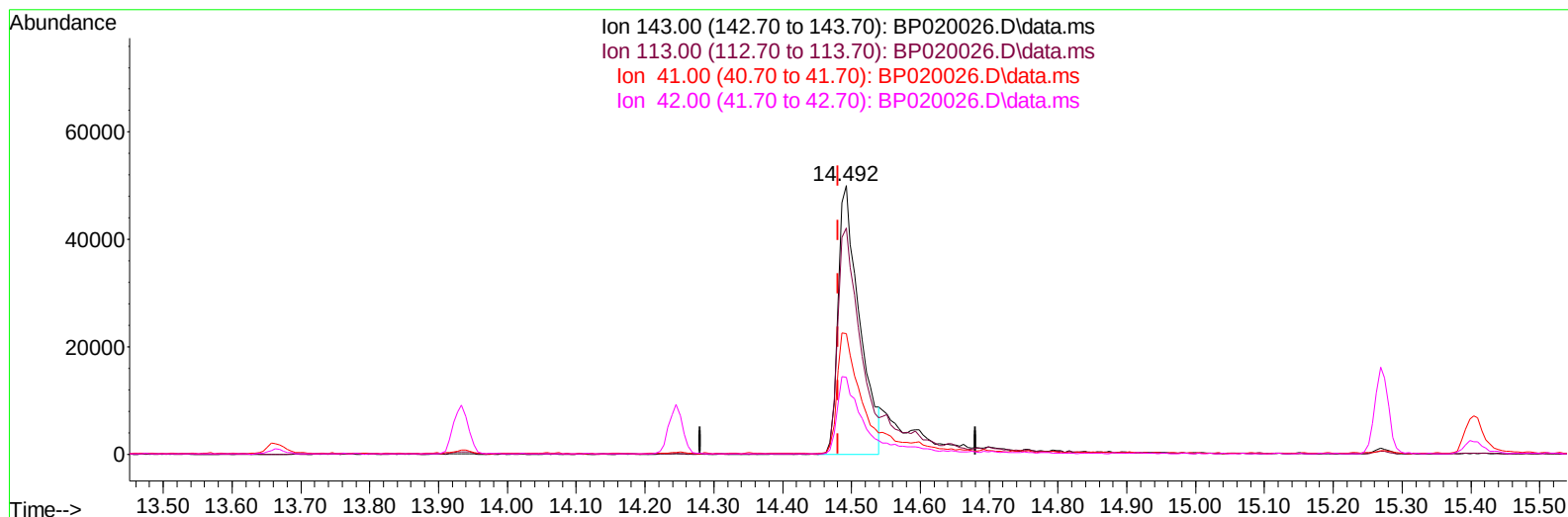
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TIC: BP020026.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.011) 23.87 ng/ul

response 107155

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	84.41
41.00	40.90	44.94
42.00	28.50	28.75

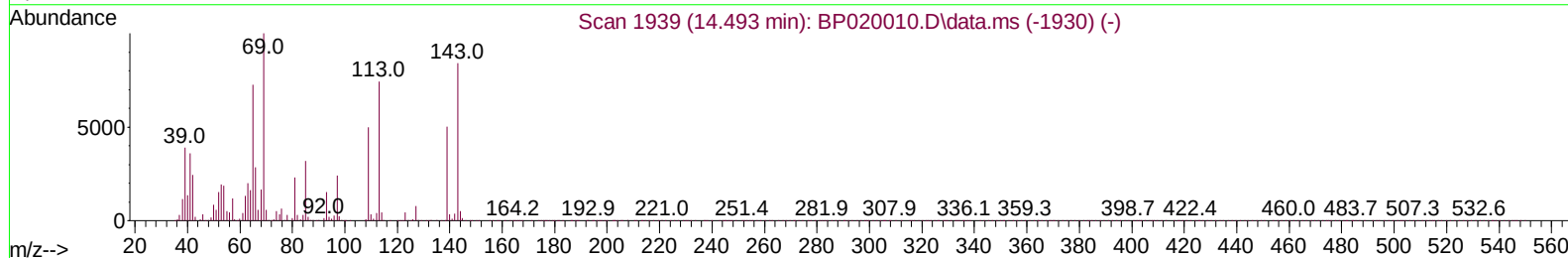
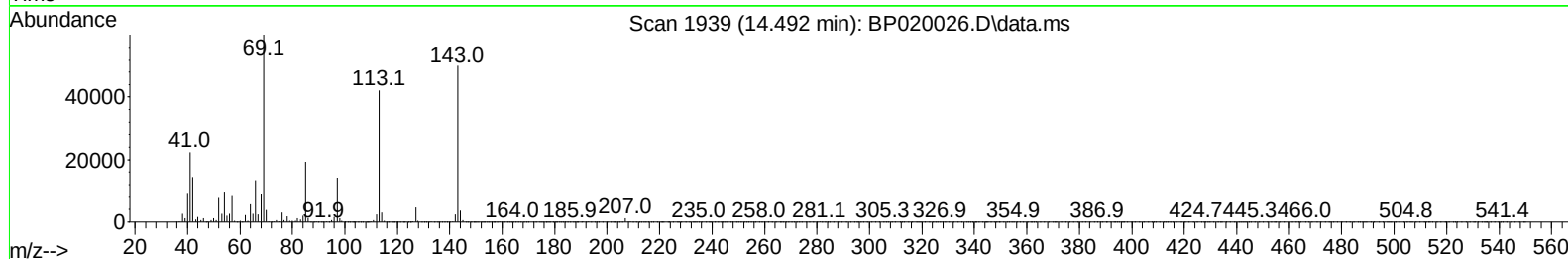
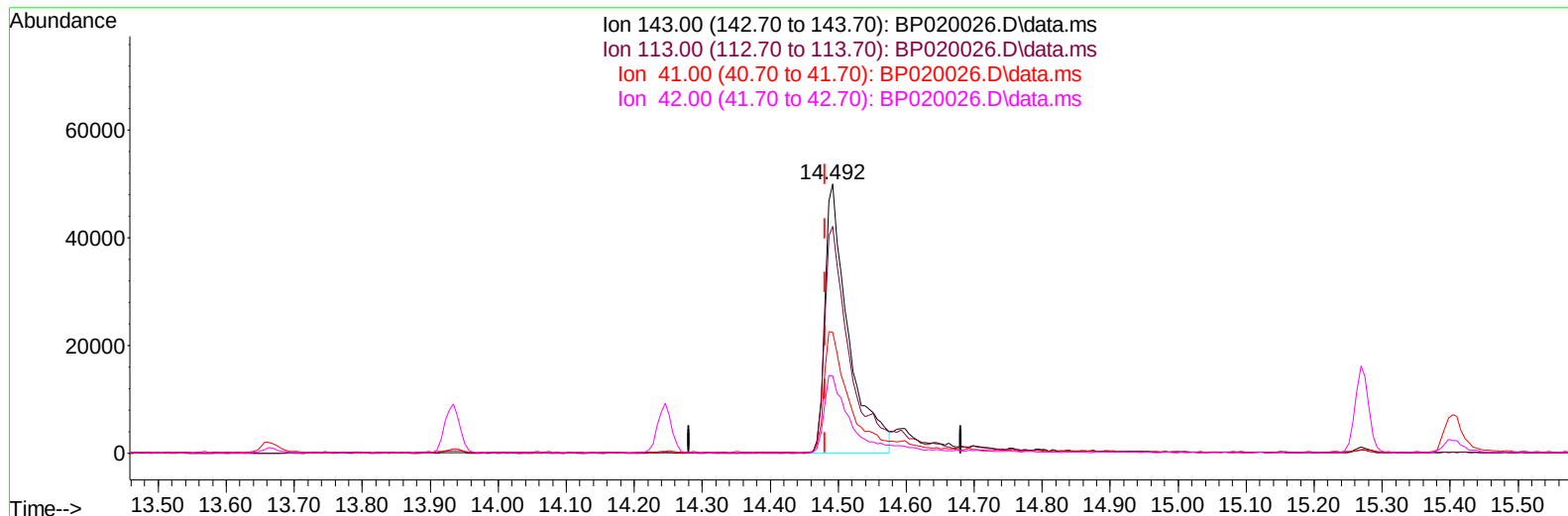
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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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TIC: BP020026.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.011) 26.72 ng/ul m

response 119962

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	84.41
41.00	40.90	44.94
42.00	28.50	28.75

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
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Quant Time: Mar 29 04:42:19 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	139159	20.000	ng/ul	0.00
20) Naphthalene-d8	10.357	136	588521	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.245	164	407294	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	835863	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	626401	20.000	ng/ul	-0.01
88) Perylene-d12	24.845	264	611705	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	23080	7.267	ng/uL	0.00
4) Pyridine-d5	3.640	84	214906	22.409	ng/ul	0.00
7) Phenol-d5	6.804	99	384390	32.752	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	272390	35.623	ng/ul	0.00
11) 2-Chlorophenol-d4	7.157	132	290010	34.407	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	296282	32.629	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	152291	33.761	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	172901	34.258	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.998	165	287463	31.187	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	14080m	1.172	ng/ul	0.05
46) Dimethylphthalate-d6	13.663	166	1047299	36.675	ng/ul	0.01
49) Acenaphthylene-d8	13.934	160	1118747	33.537	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	119962m	26.722	ng/ul	0.01
60) Fluorene-d10	15.275	176	858799	35.304	ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.404	200	155209	31.021	ng/ul	-0.02
73) Anthracene-d10	17.180	188	1283708	34.694	ng/ul	0.01
81) Pyrene-d10	19.574	212	1397148	38.016	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.609	264	990676	33.711	ng/ul	0.00

Target Compounds Qvalue

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

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