

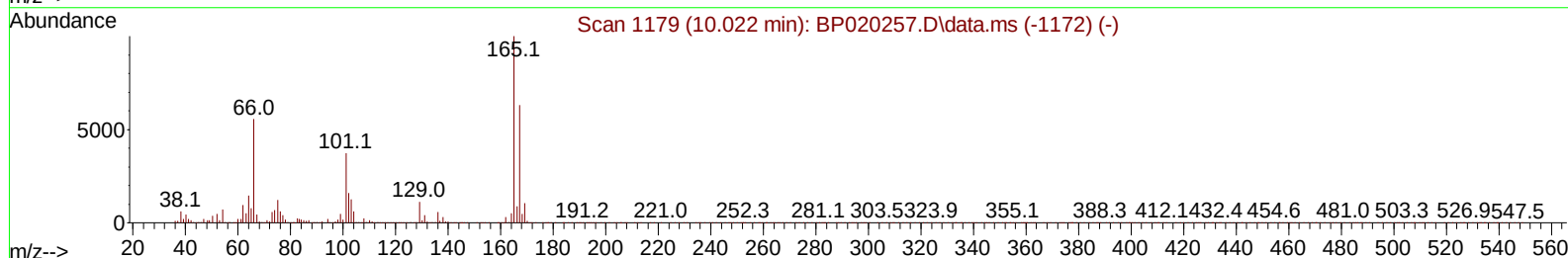
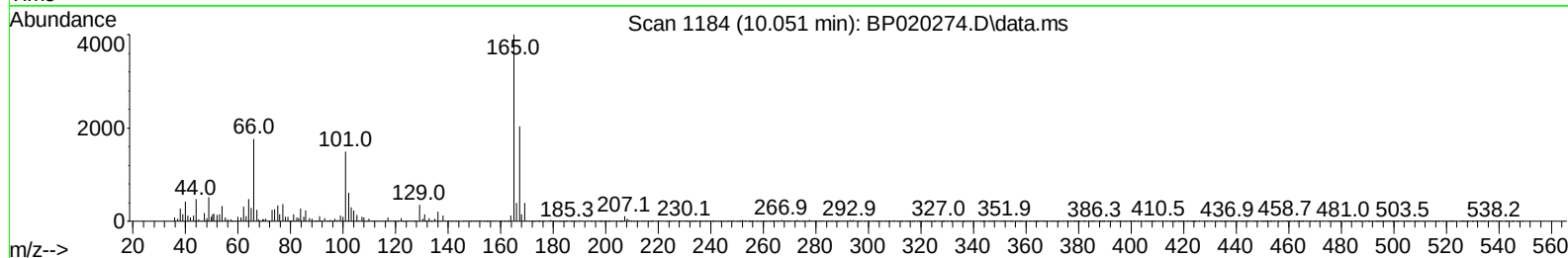
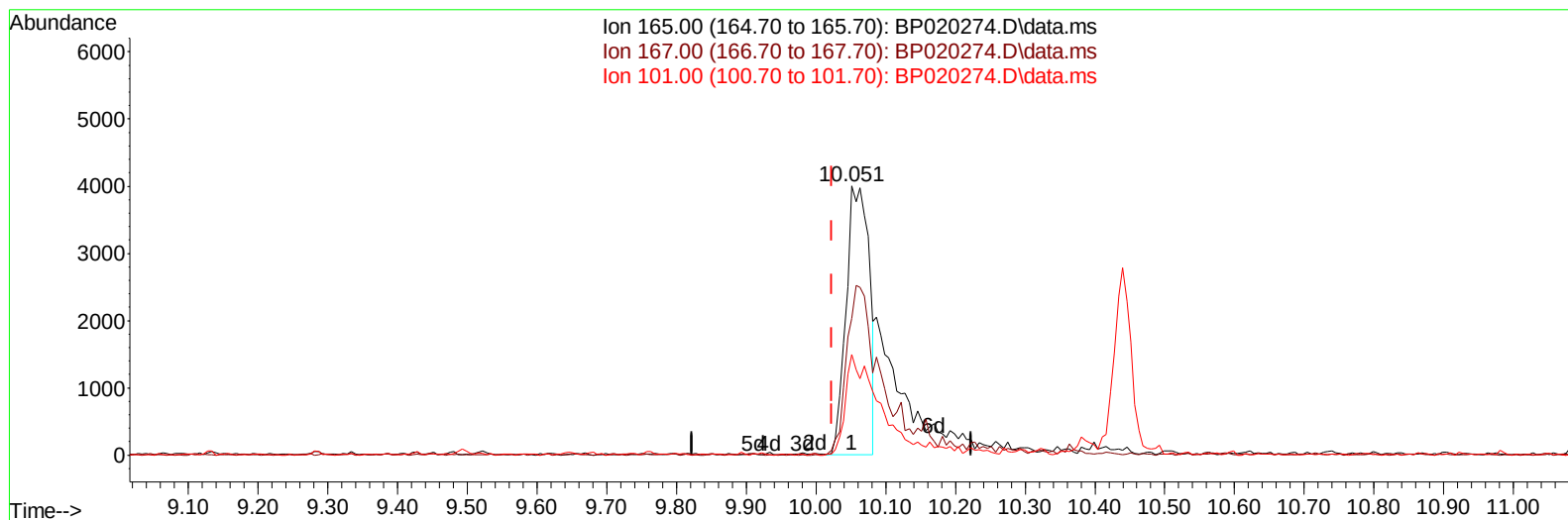
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020274.D
 Acq On : 09 May 2024 19:40
 Operator : MA/JU
 Sample : P2369-DL 10X
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43M6DL

Manual IntegrationsAPPROVED

Quant Time: May 09 22:45:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/10/2024
 Supervised By :mohammad ahmed 05/11/2024



TIC: BP020274.D\data.ms

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

10.051min (+ 0.029) 1.74 ng/ul

response 9111

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	63.00	50.89
101.00	40.90	37.29
0.00	0.00	0.00

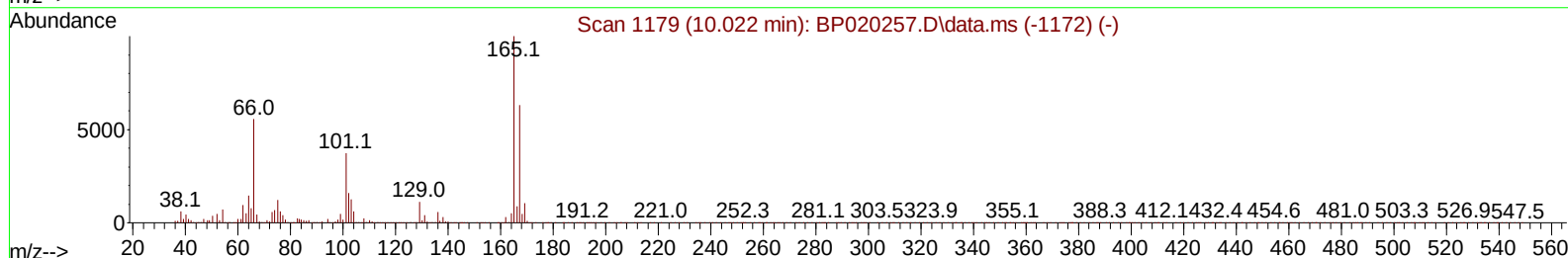
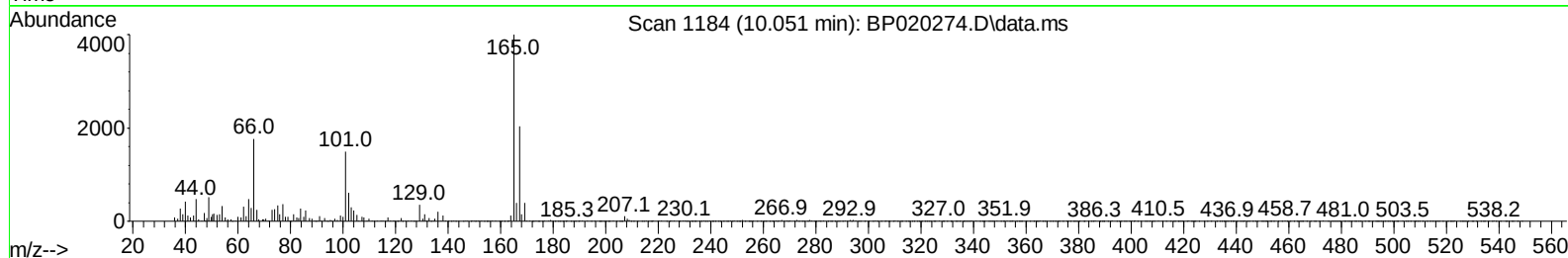
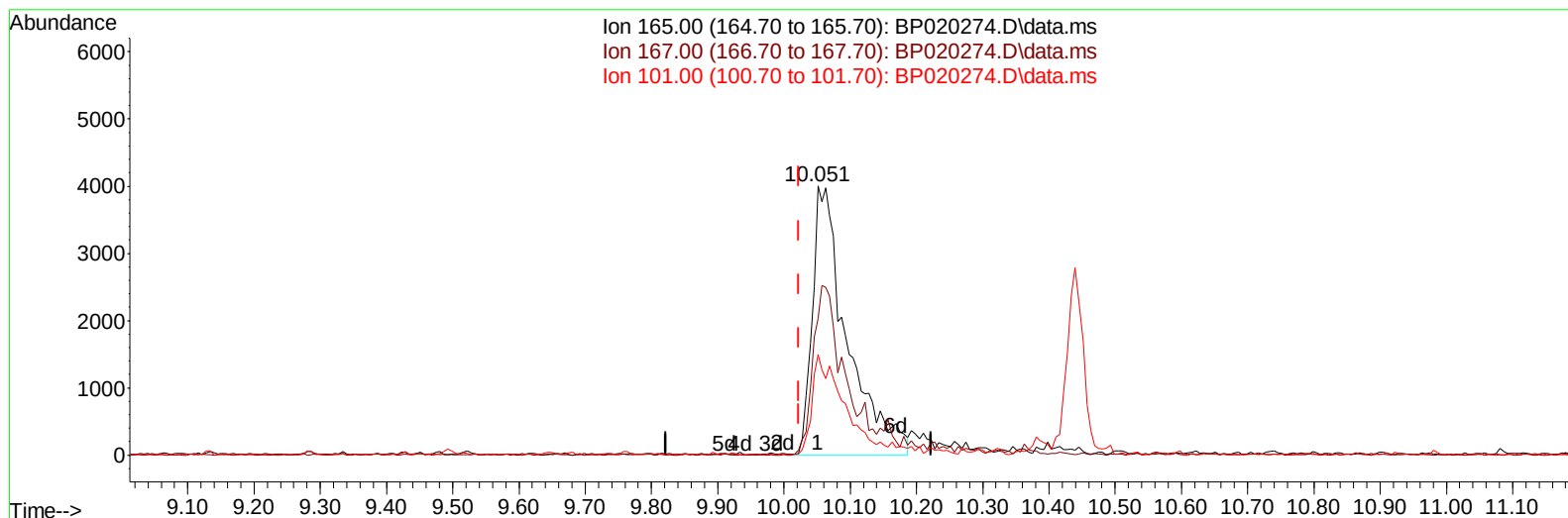
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TIC: BP020274.D\data.ms

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

10.051min (+ 0.029) 2.78 ng/ul m

response 14579

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	63.00	50.89
101.00	40.90	37.29
0.00	0.00	0.00

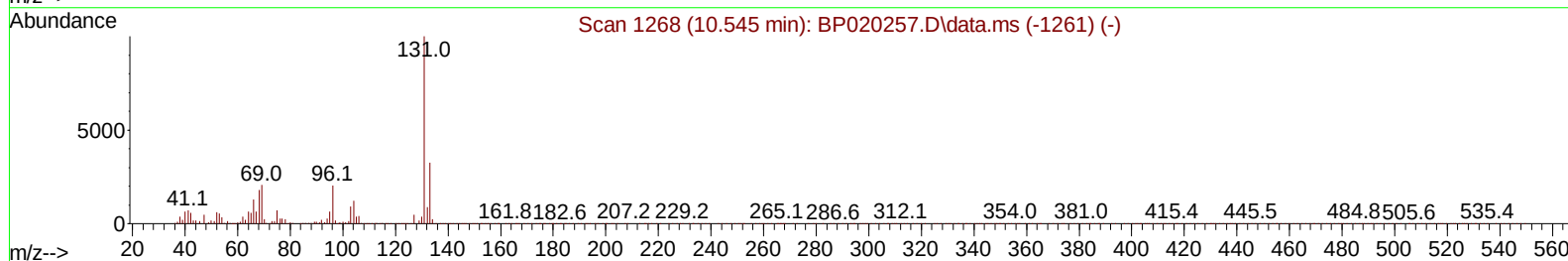
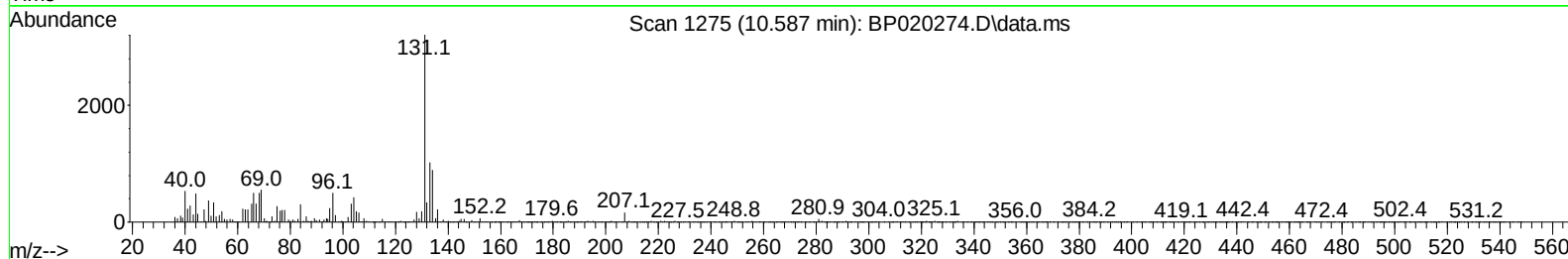
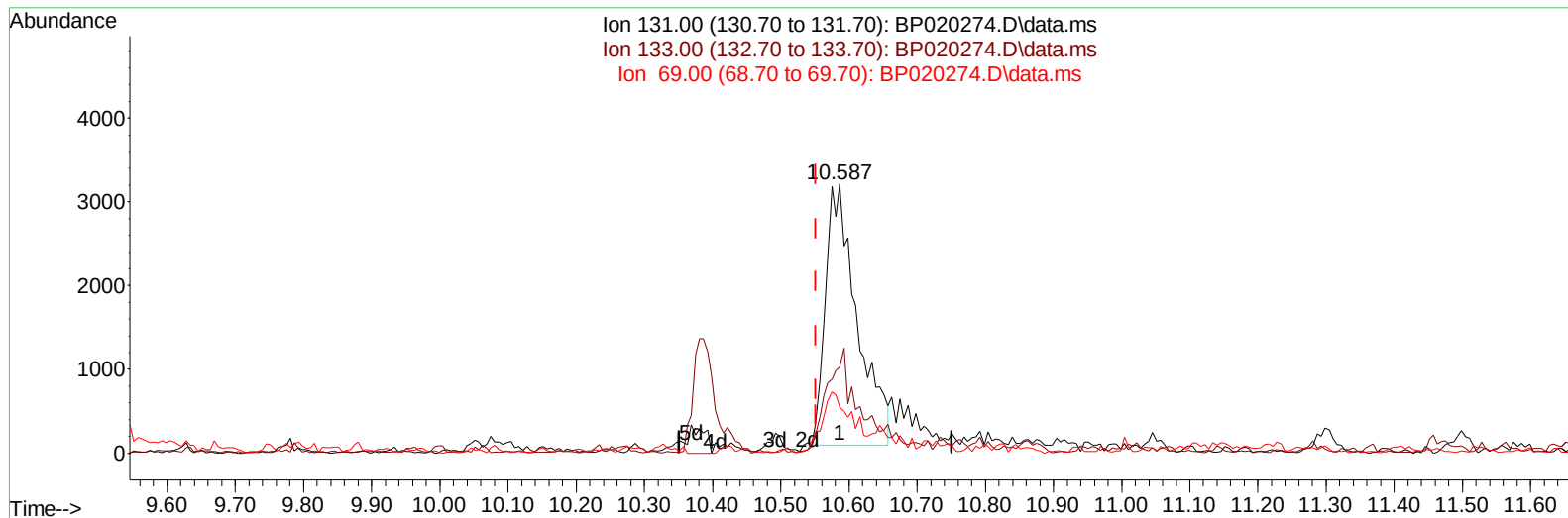
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TIC: BP020274.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.587min (+ 0.035) 1.36 ng/ul

response 10000

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.40	32.01
69.00	20.50	17.16
0.00	0.00	0.00

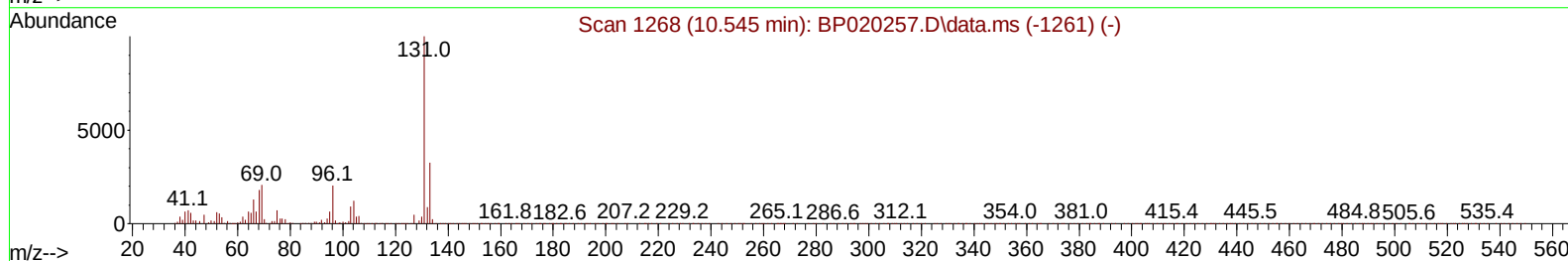
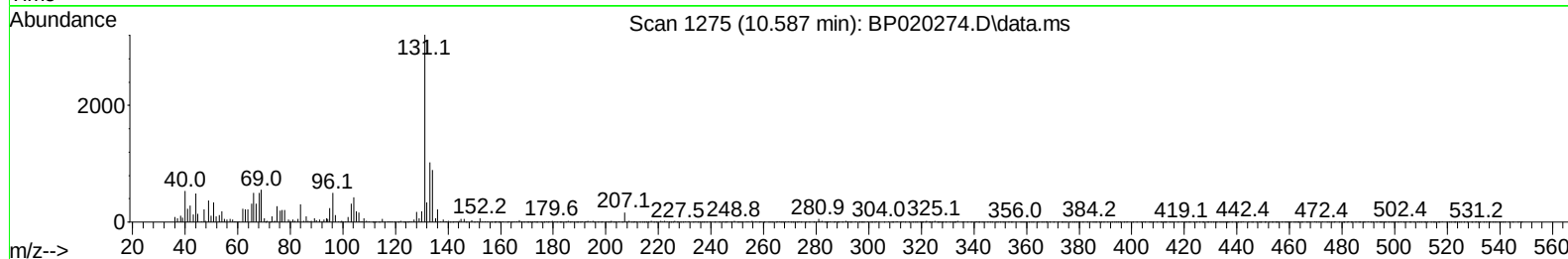
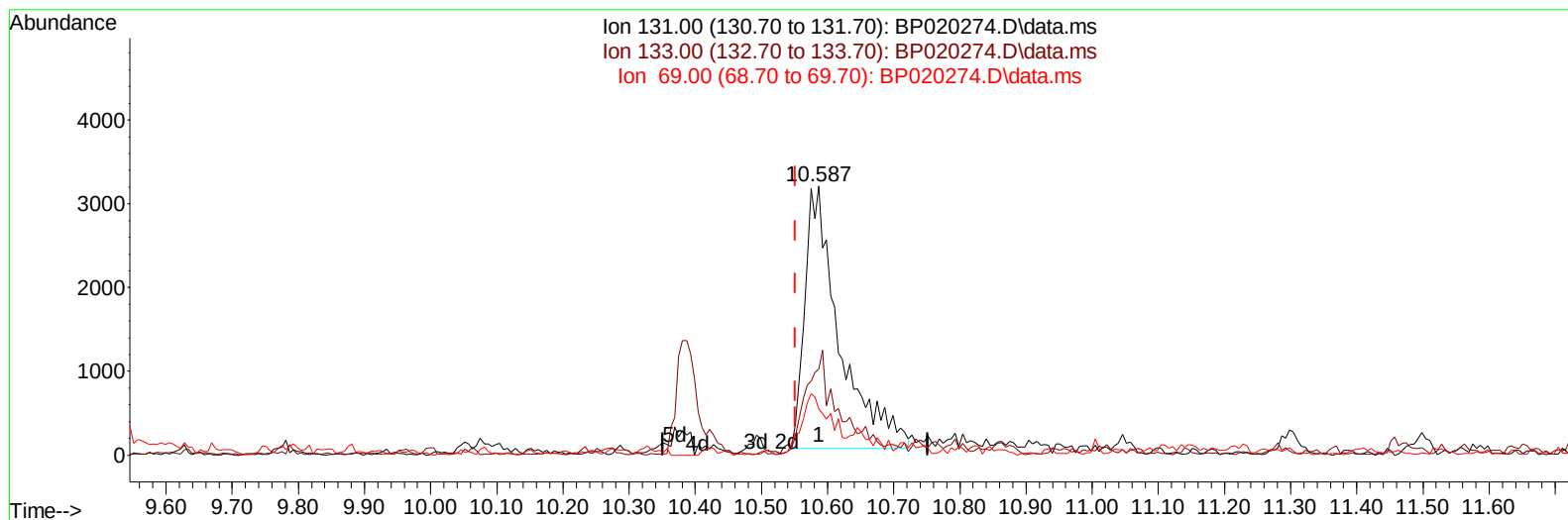
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Misc :
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Instrument :
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TIC: BP020274.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.587min (+ 0.035) 1.56 ng/ul m

response 11443

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.40	32.01
69.00	20.50	17.16
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/10/2024
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Quant Time: May 09 22:58:23 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	77560	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	335444	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.292	164	237022	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.133	188	522276	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	533138	20.000	ng/ul	0.00
88) Perylene-d12	24.992	264	620377	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	879	0.479	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.828	99	14514	2.179	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.981	67	8118	2.006	ng/ul	0.01
11) 2-Chlorophenol-d4	7.163	132	12354	2.594	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	11302	2.099	ng/ul	0.02
21) Nitrobenzene-d5	8.793	128	5693	2.266	ng/ul	0.01
24) 2-Nitrophenol-d4	9.498	143	6612	2.299	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.051	165	14579m	2.777	ng/ul	0.03
31) 4-Chloroaniline-d4	10.587	131	11443m	1.558	ng/ul	0.04
46) Dimethylphthalate-d6	13.716	166	55021	3.179	ng/ul	0.00
49) Acenaphthylene-d8	13.986	160	56103	2.934	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.322	176	48872	3.385	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	4905	1.480	ng/ul	0.01
73) Anthracene-d10	17.239	188	79462	3.517	ng/ul	0.00
81) Pyrene-d10	19.645	212	97427	3.126	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.768	264	100418	3.348	ng/ul	0.00

Target Compounds Qvalue

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

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