

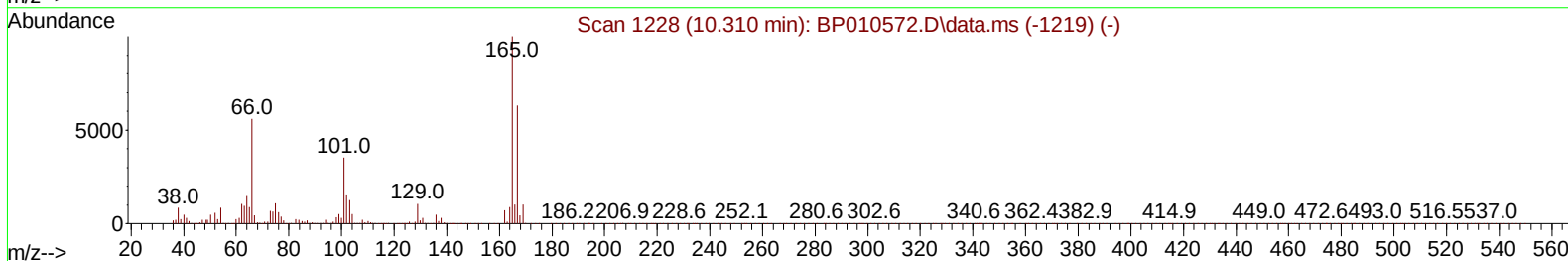
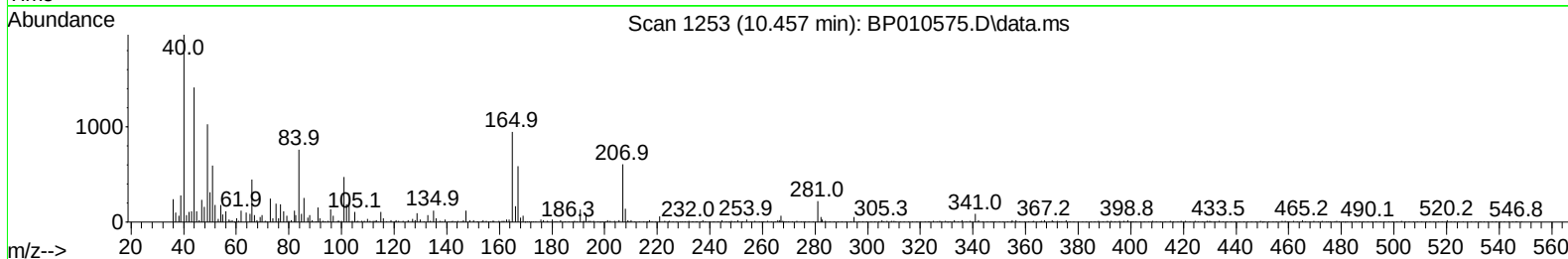
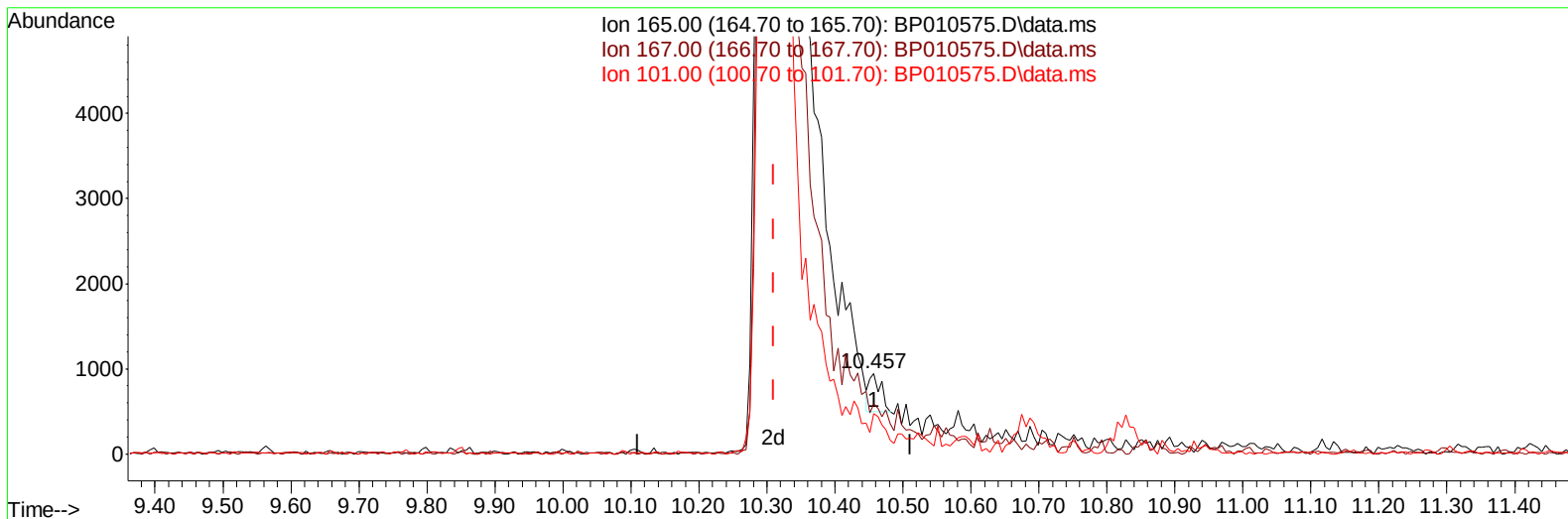
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052722\
Data File : BP010575.D
Acq On : 27 May 2022 11:50
Operator : CG/JU
Sample : N2883-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4K4

Manual IntegrationsAPPROVED

Quant Time: May 28 00:14:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 24 05:50:40 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/31/2022
Supervised By :mohammad ahmed 05/31/2022



TIC: BP010575.D\data.ms

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

10.457min (+ 0.147) 0.05 ng/u1

response 527

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	61.77
101.00	49.80	49.84
0.00	0.00	0.00

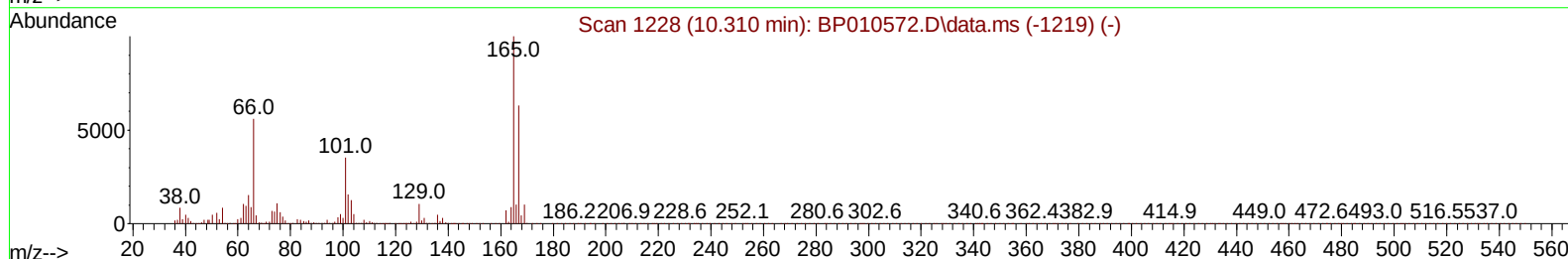
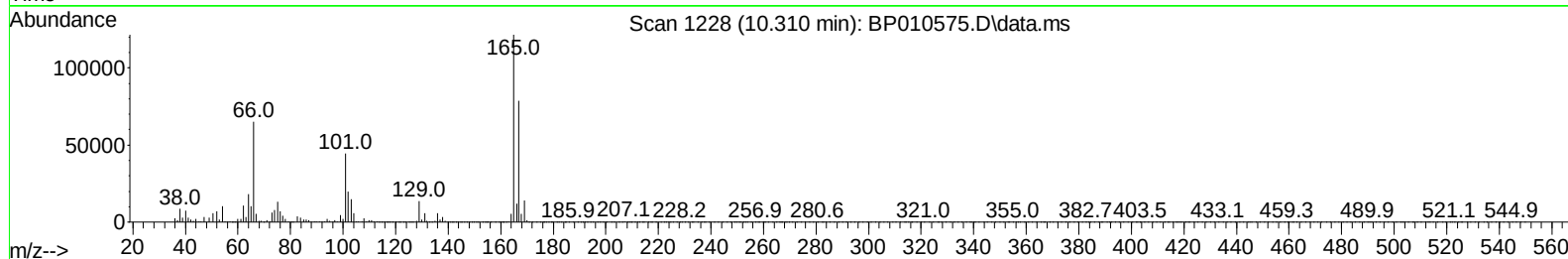
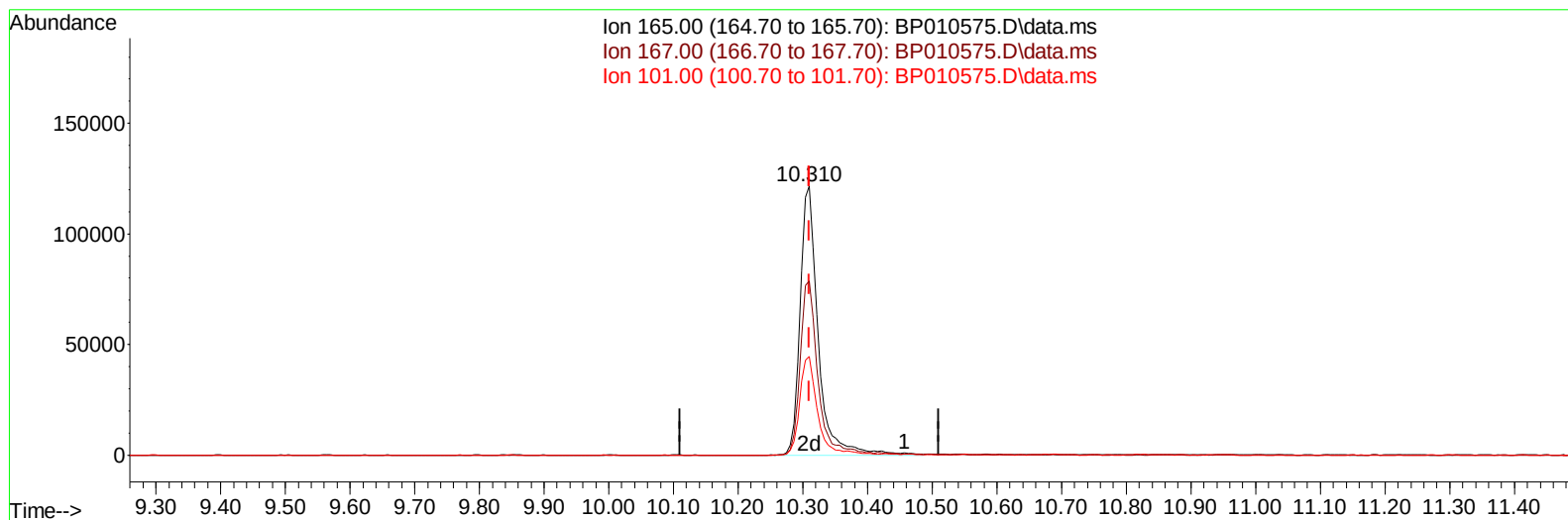
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TIC: BP010575.D\data.ms

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

10.310min (-0.000) 23.08 ng/ul m

response 236656

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	64.77
101.00	49.80	36.60#
0.00	0.00	0.00

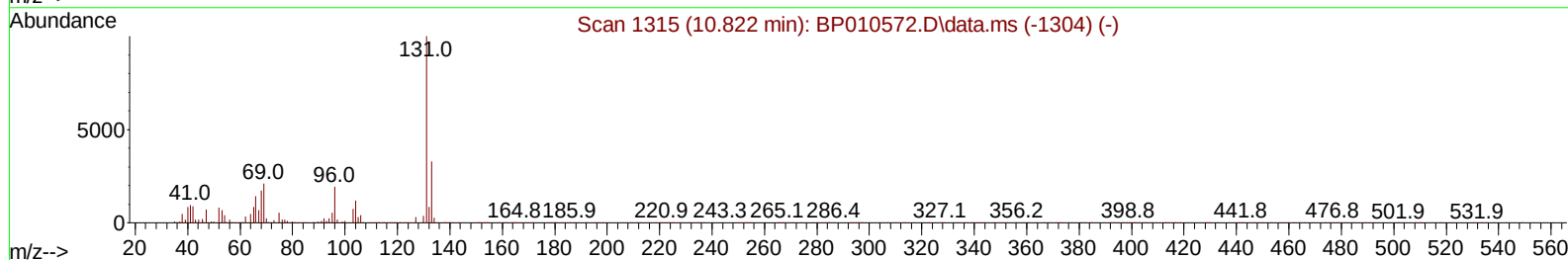
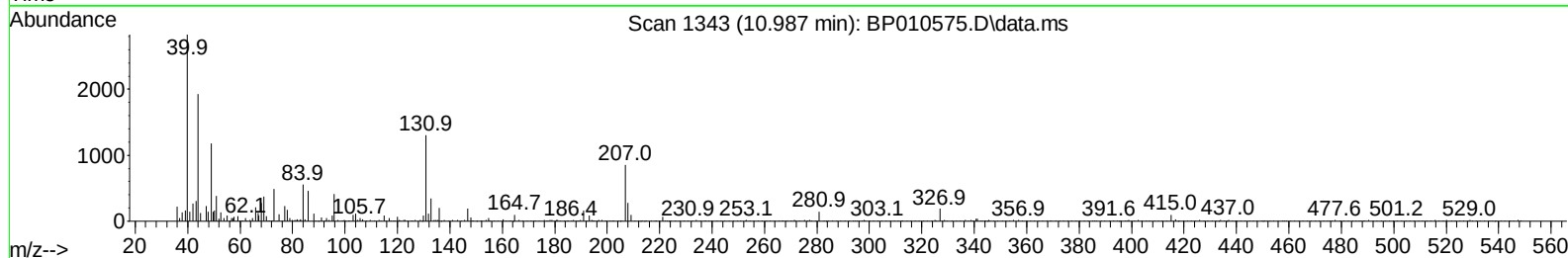
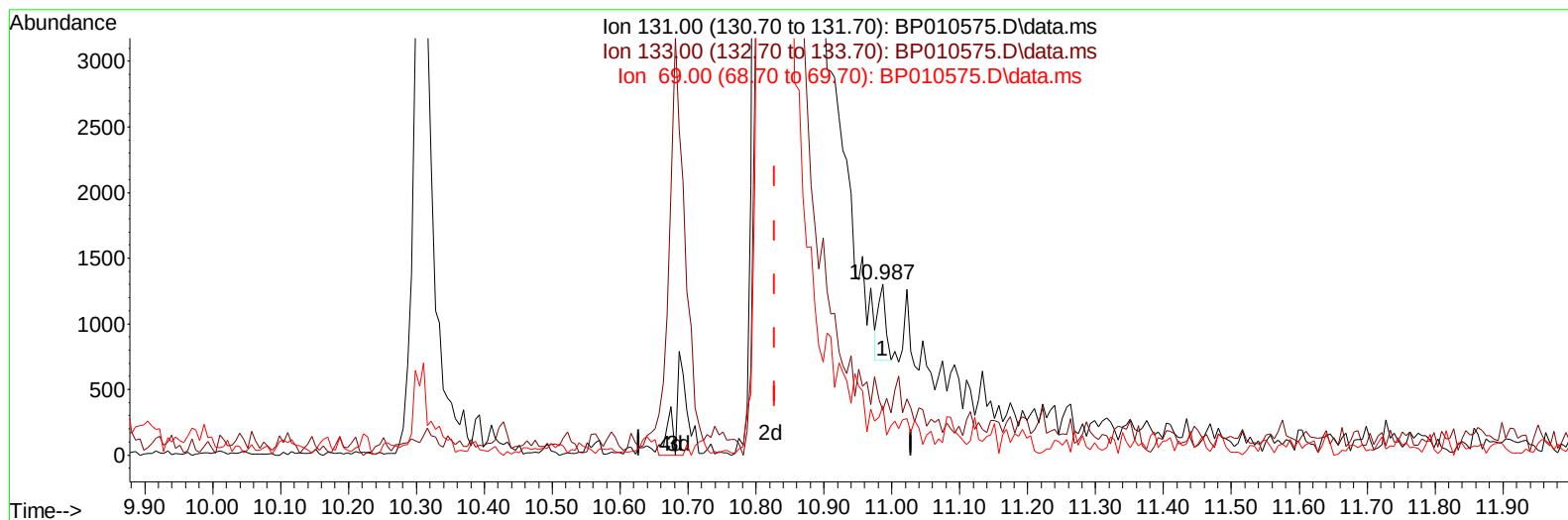
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Instrument :
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TIC: BP010575.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.987min (+ 0.159) 0.03 ng/ul

response 429

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	26.34
69.00	26.10	28.80
0.00	0.00	0.00

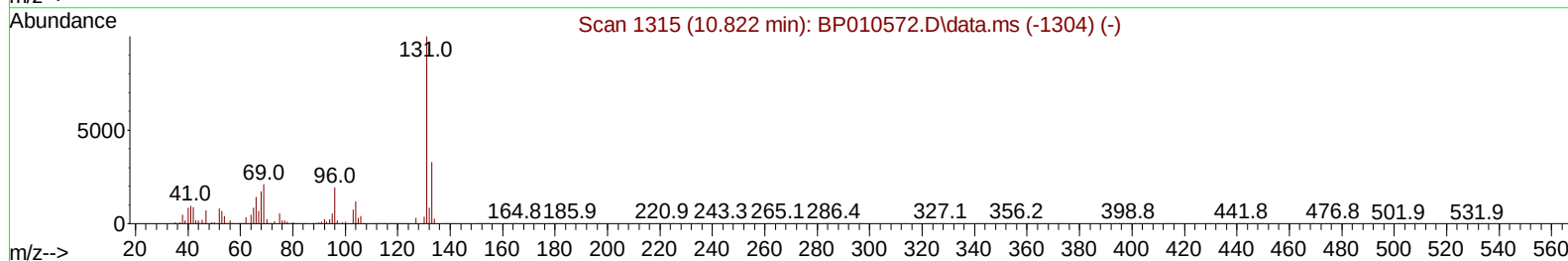
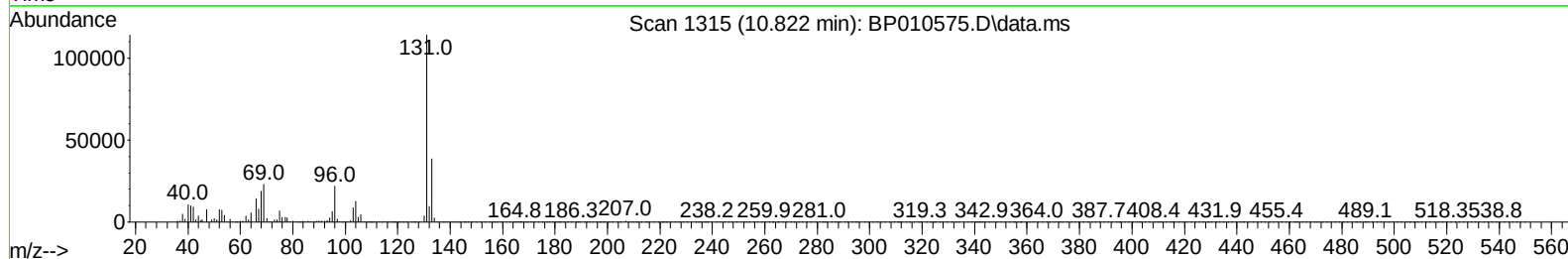
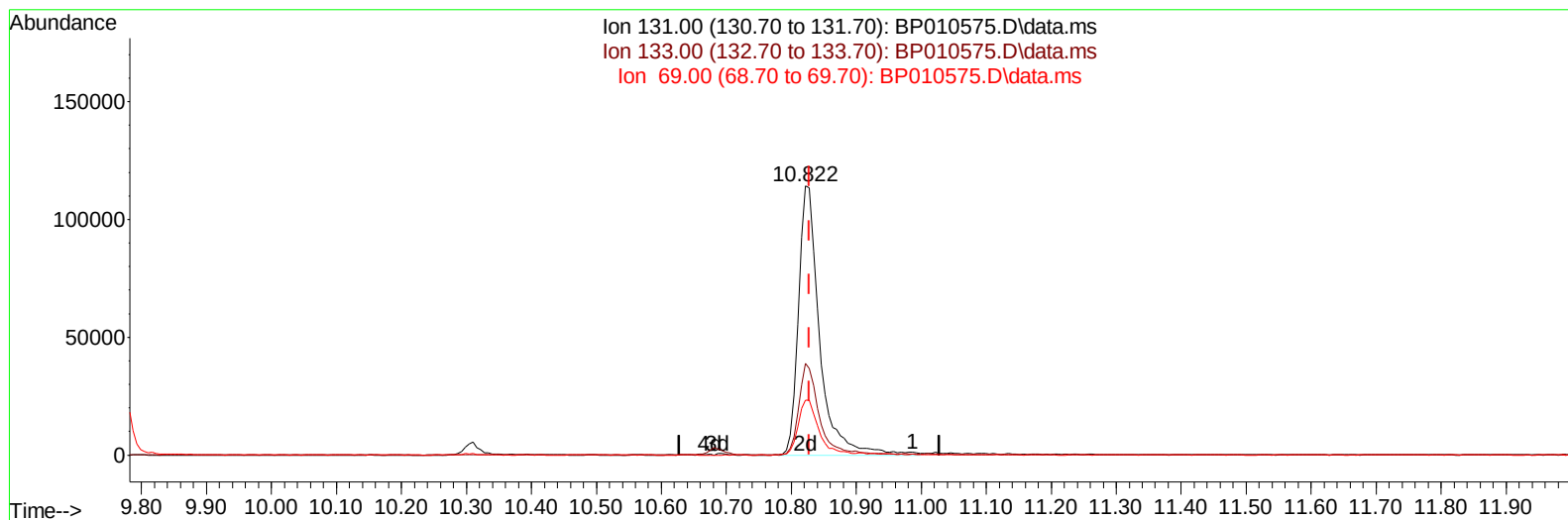
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Instrument :
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TIC: BP010575.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.822min (-0.006) 16.71 ng/ul m

response 254841

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	34.04
69.00	26.10	20.30#
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	7.881	152	150131	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	661224	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.516	164	462237	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	1019270	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	836441	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	666725	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	15907	4.516	ng/uL	0.00
4) Pyridine-d5	3.752	84	221222	22.120	ng/ul	0.00
7) Phenol-d5	7.052	99	284675	22.592	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	213364	25.202	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	236821	24.561	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	246918	23.058	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	127191	24.854	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	134874	24.870	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	236656m	23.078	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	254841m	16.714	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	908648	26.681	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	1012716	25.780	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	84468	13.935	ng/ul	0.00
60) Fluorene-d10	15.510	176	771029	26.023	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	93687	15.716	ng/ul	0.00
73) Anthracene-d10	17.369	188	1191118	25.807	ng/ul	0.00
81) Pyrene-d10	19.604	212	1385193	31.183	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	887323	26.504	ng/ul	0.00

Target Compounds Qvalue

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

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