

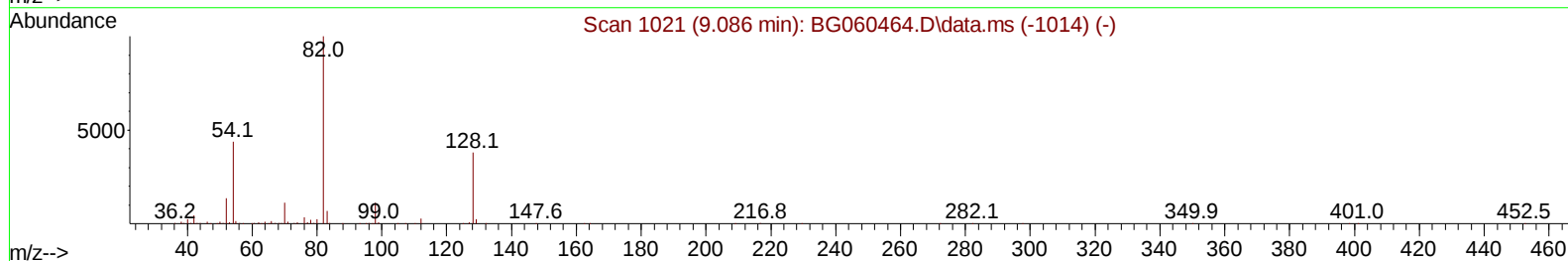
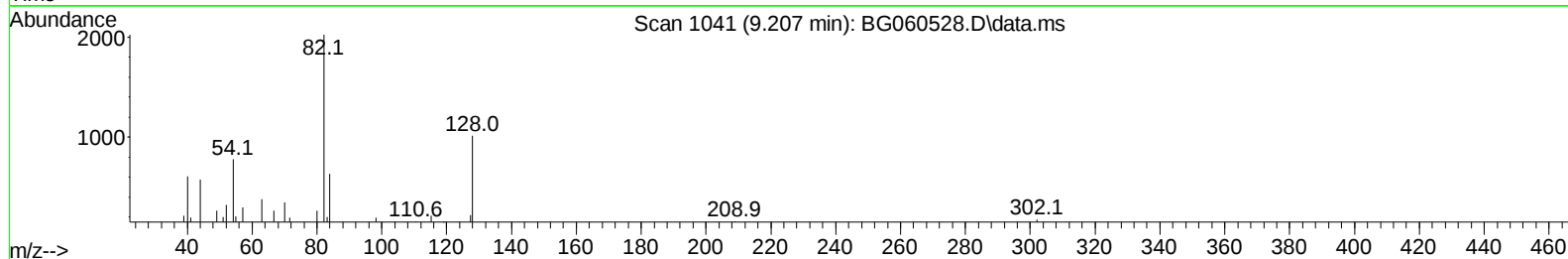
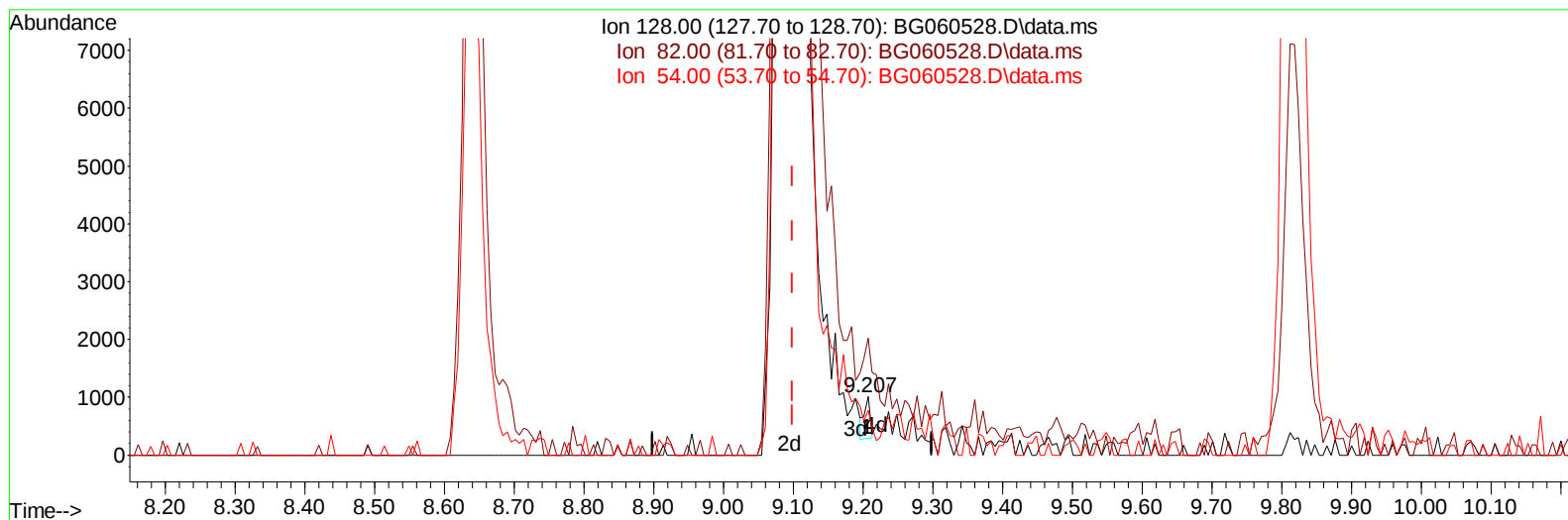
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020124\
Data File : BG060528.D
Acq On : 1 Feb 2024 4:40
Operator : MA/JU
Sample : P1289-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF9

Manual IntegrationsAPPROVED

Quant Time: Feb 01 05:15:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 31 22:04:13 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/01/2024
Supervised By :Sohil Jodhani 02/02/2024



TIC: BG060528.D\data.ms

(21) Nitrobenzene-d5 (S)

9.207min (+ 0.108) 0.10 ng/u1

response 417

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	238.70	199.51
54.00	95.30	76.73
0.00	0.00	0.00

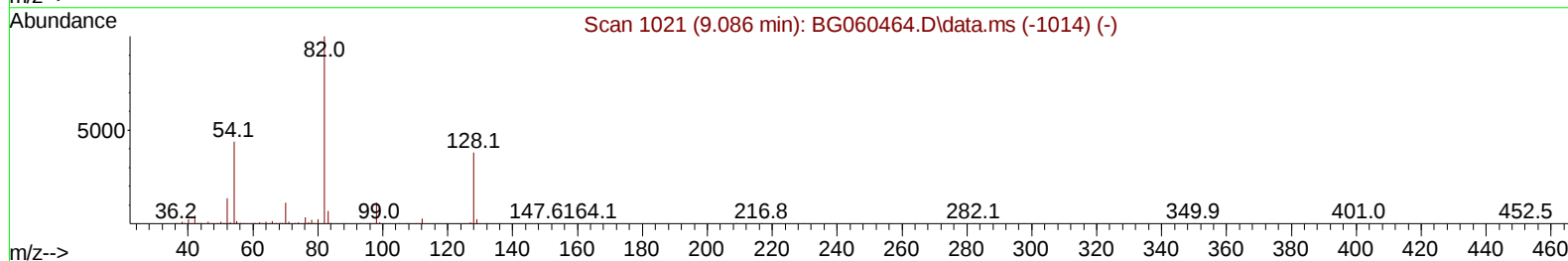
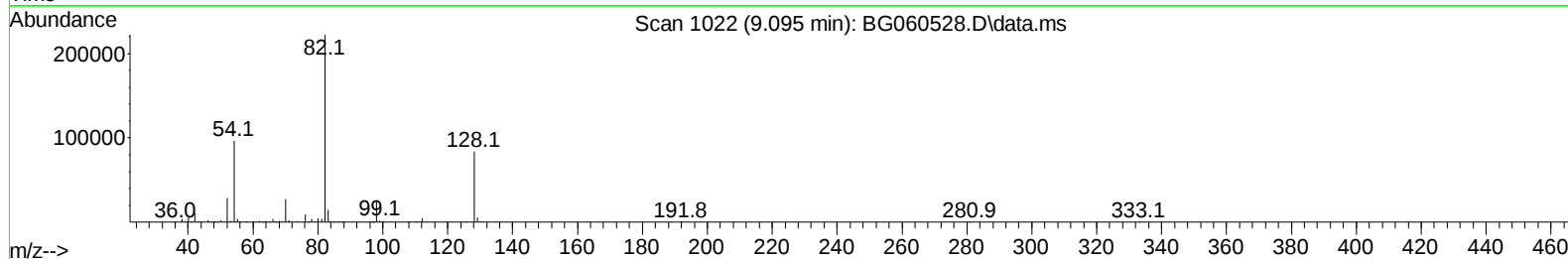
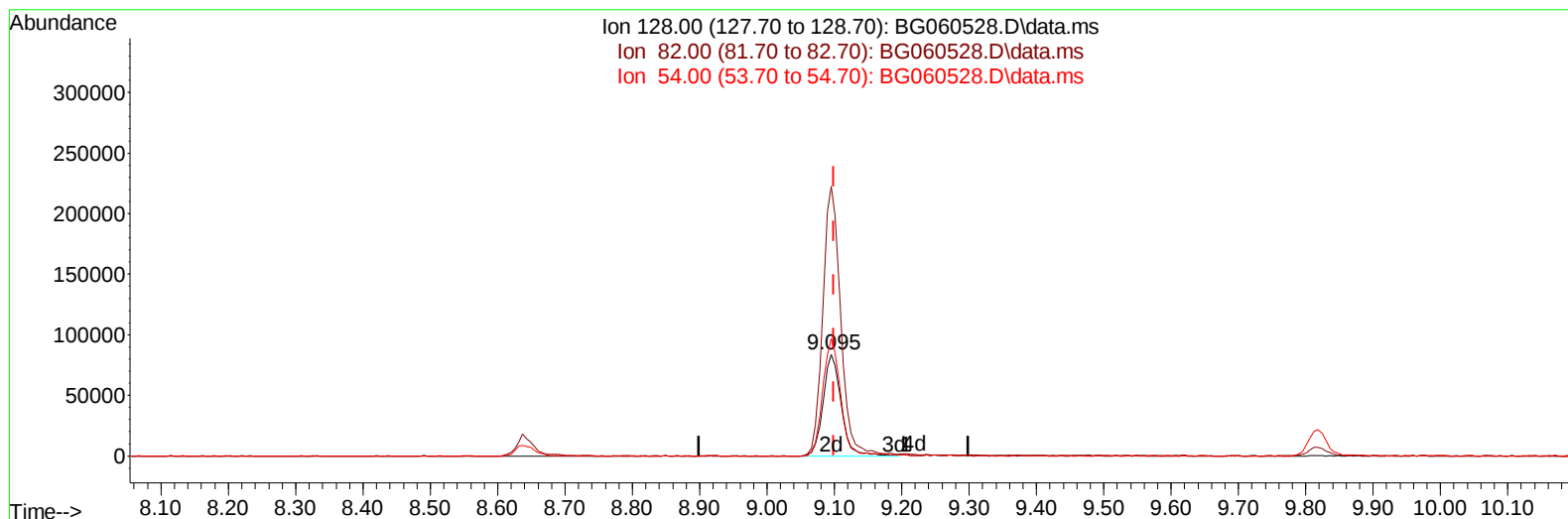
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TIC: BG060528.D\data.ms

(21) Nitrobenzene-d5 (S)

9.095min (-0.004) 37.78 ng/ul m

response 156445

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	238.70	265.70
54.00	95.30	115.81#
0.00	0.00	0.00

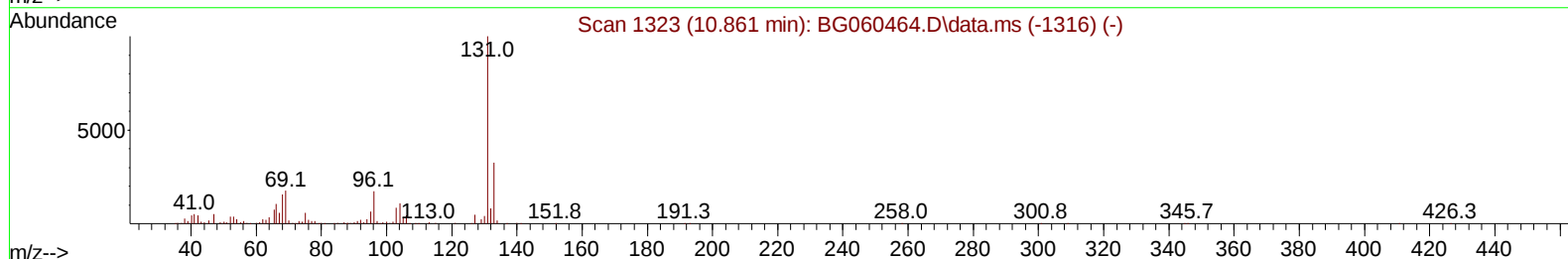
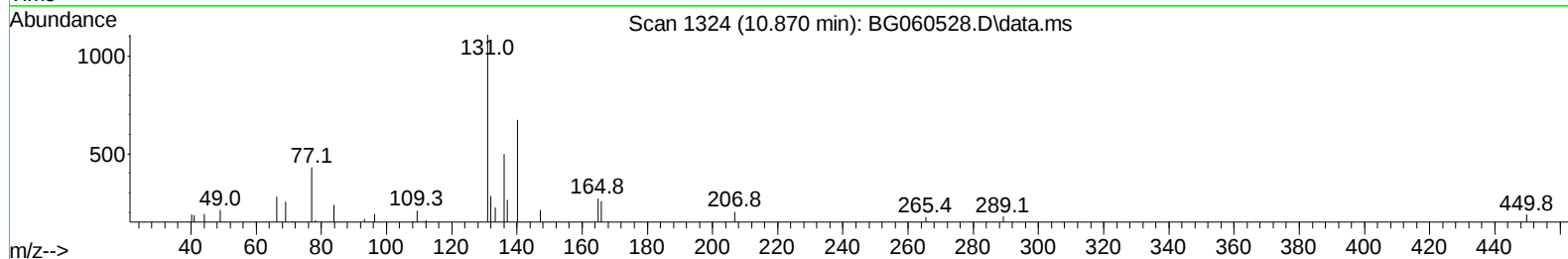
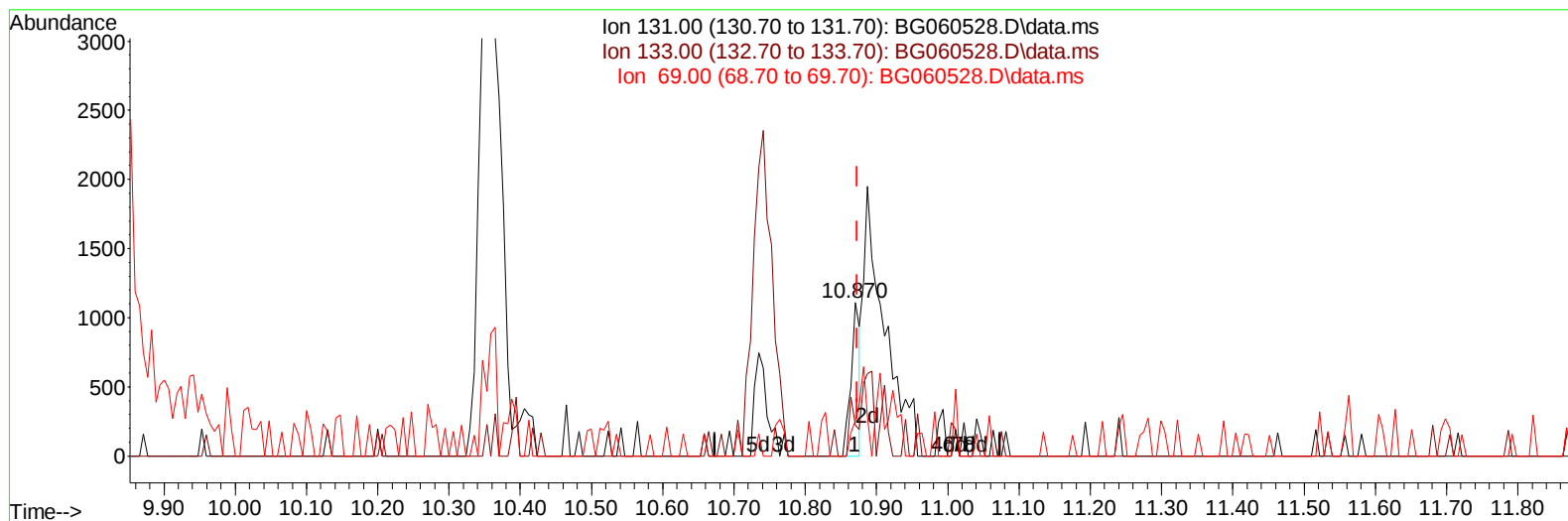
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG060528.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.870min (-0.004) 0.07 ng/u1

response 987

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.60	20.29#
69.00	19.60	22.90
0.00	0.00	0.00

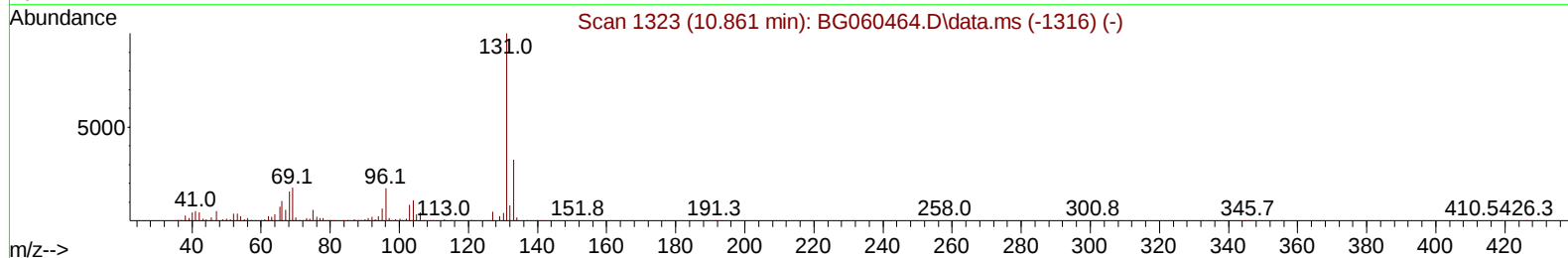
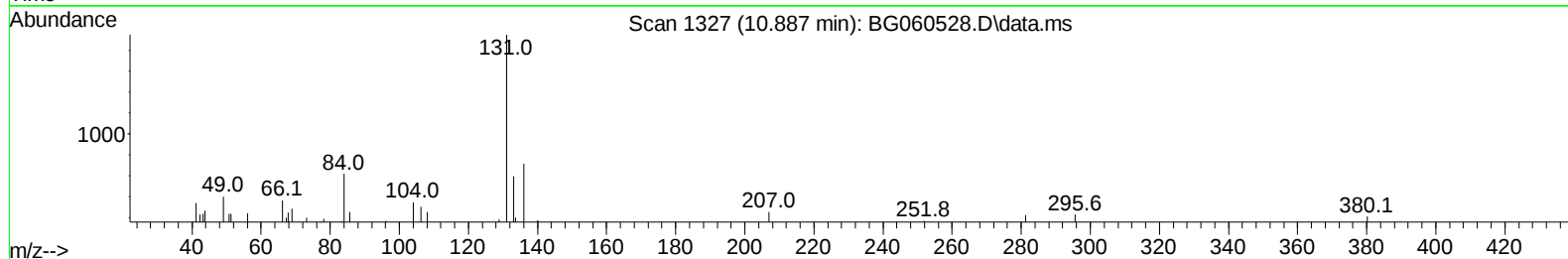
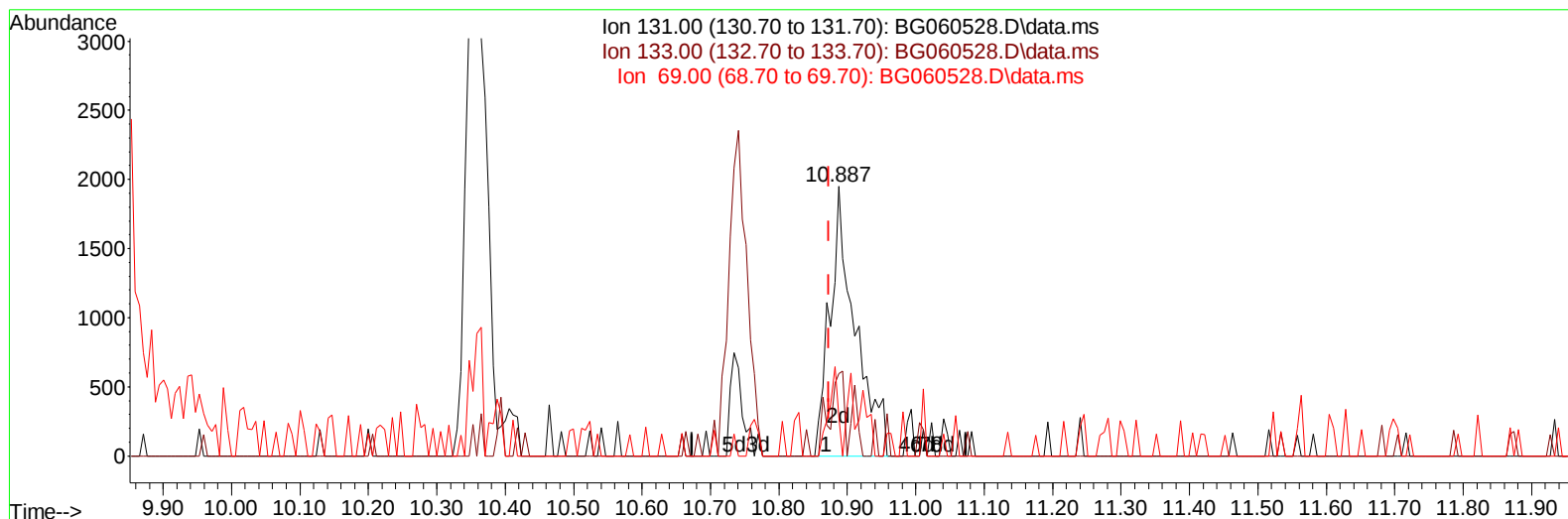
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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TIC: BG060528.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.887min (+ 0.014) 0.38 ng/ul m

response 4994

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.60	30.48
69.00	19.60	14.57#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.938	152	128086	20.000	ng/ul	0.00
20) Naphthalene-d8	10.735	136	561275	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.571	164	502118	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	1366499	20.000	ng/ul	0.00
79) Chrysene-d12	21.581	240	1362087	20.000	ng/ul	0.00
88) Perylene-d12	24.754	264	1510389	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.373	96	10879	4.256	ng/uL	0.00
4) Pyridine-d5	3.796	84	5292	0.654	ng/ul	0.01
7) Phenol-d5	7.098	99	73026	6.003	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.262	67	244085	31.930	ng/ul	0.00
11) 2-Chlorophenol-d4	7.468	132	214887	26.453	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	155518	15.520	ng/ul	0.00
21) Nitrobenzene-d5	9.095	128	156445m	37.785	ng/ul	0.00
24) 2-Nitrophenol-d4	9.818	143	187509	38.763	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.359	165	351671	34.957	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	4994m	0.378	ng/ul	0.01
46) Dimethylphthalate-d6	13.972	166	1537895	37.430	ng/ul	0.00
49) Acenaphthylene-d8	14.266	160	1475766	32.020	ng/ul	0.00
54) 4-Nitrophenol-d4	14.777	143	36372	6.181	ng/ul	0.00
60) Fluorene-d10	15.564	176	1418814	39.410	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.682	200	280087	33.101	ng/ul	0.00
73) Anthracene-d10	17.415	188	2298289	35.480	ng/ul	0.00
81) Pyrene-d10	19.707	212	2911364	35.663	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.536	264	2724955	34.392	ng/ul	0.00

Target Compounds Qvalue

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

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