

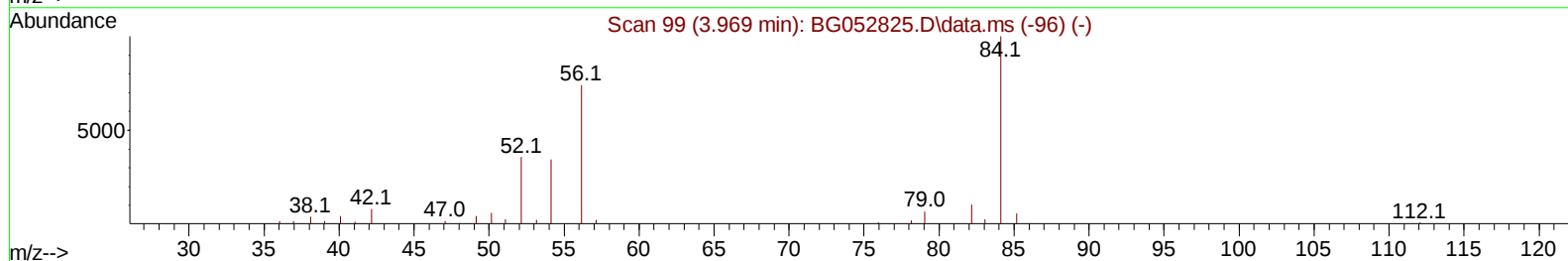
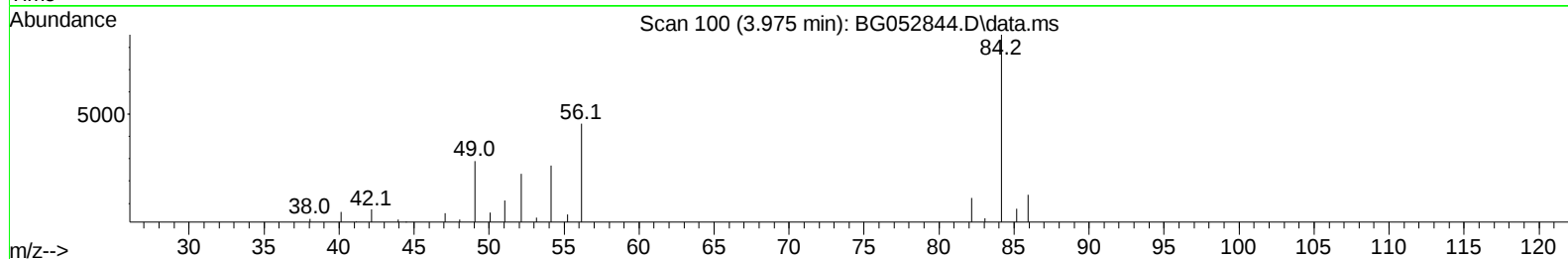
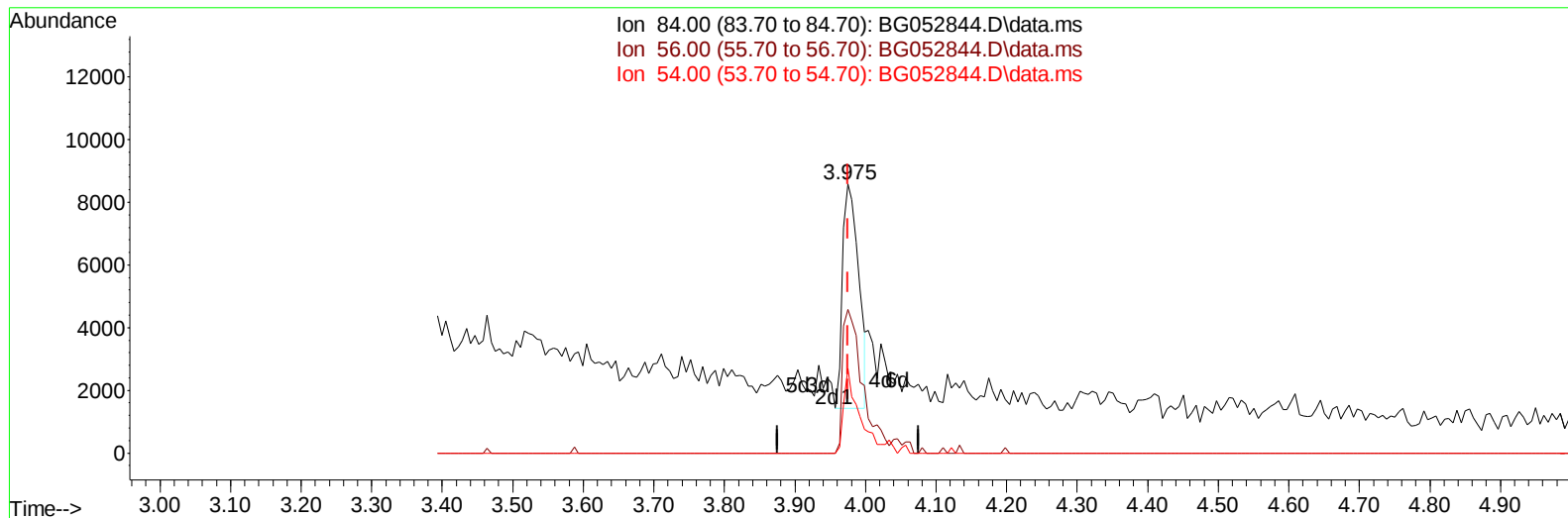
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
Data File : BG052844.D
Acq On : 26 Mar 2022 1:09
Operator : CG/JU
Sample : N2079-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW609

Manual IntegrationsAPPROVED

Quant Time: Mar 26 02:09:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 23 13:16:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/28/2022
Supervised By :mohammad ahmed 03/29/2022



TIC: BG052844.D\data.ms

(4) Pyridine-d5 (S)

3.975min (-0.001) 6.18 ng/u1

response 11431

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	53.31
54.00	28.70	31.59
0.00	0.00	0.00

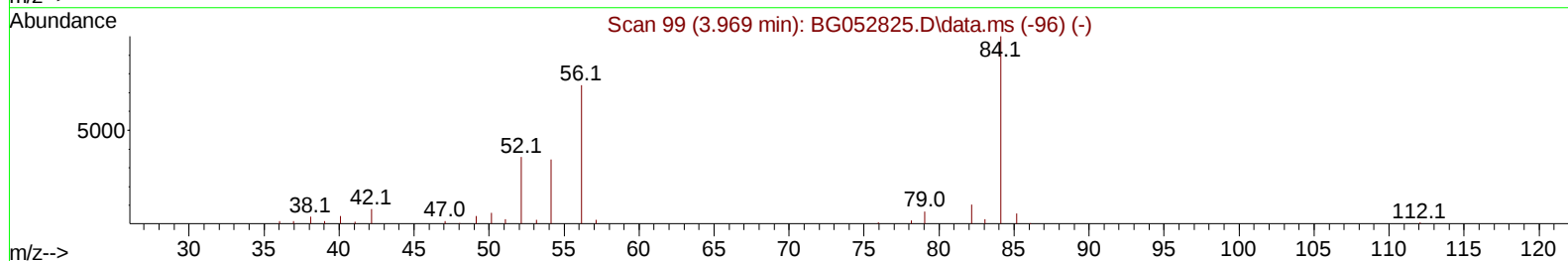
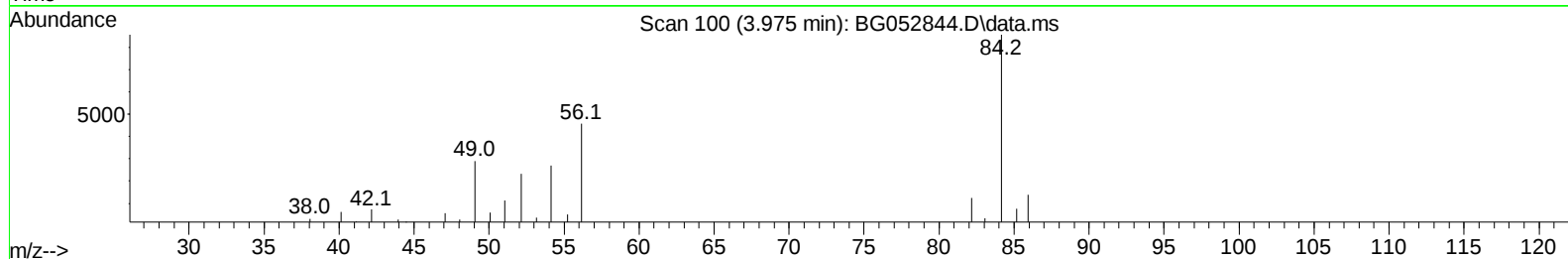
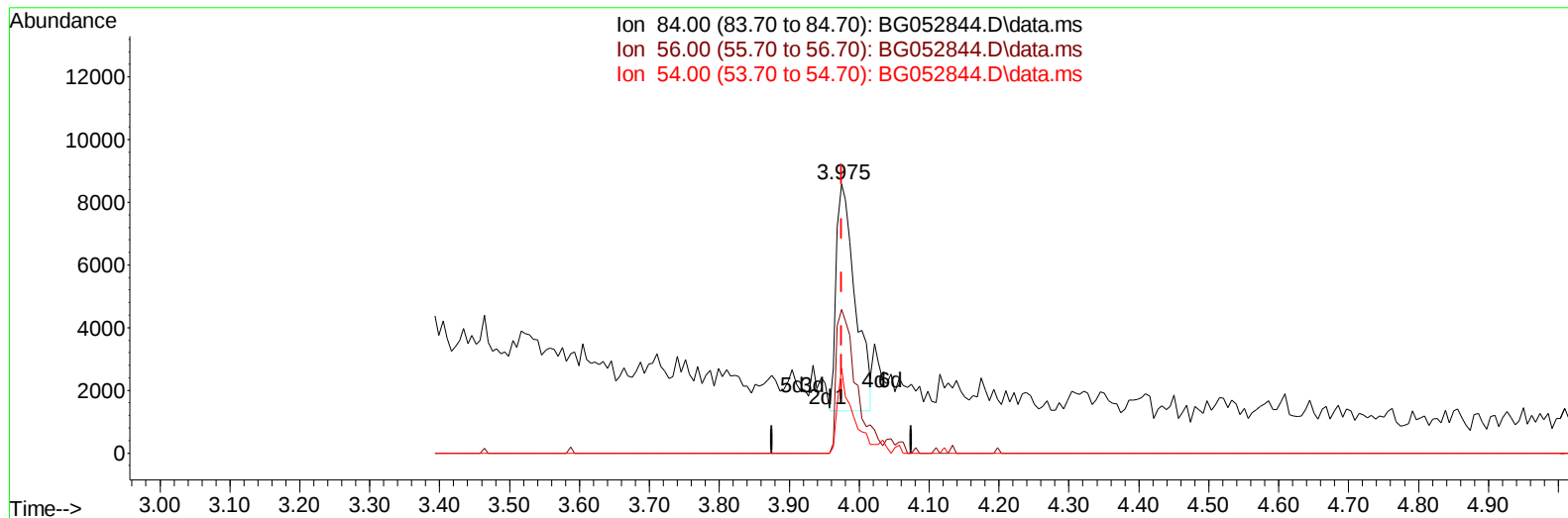
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TIC: BG052844.D\data.ms

(4) Pyridine-d5 (S)

3.975min (-0.001) 7.37 ng/ul m

response 13622

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	53.31
54.00	28.70	31.59
0.00	0.00	0.00

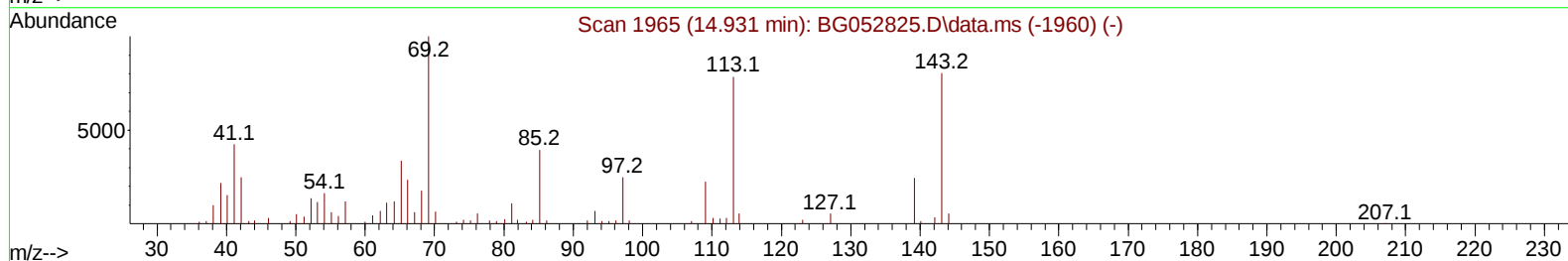
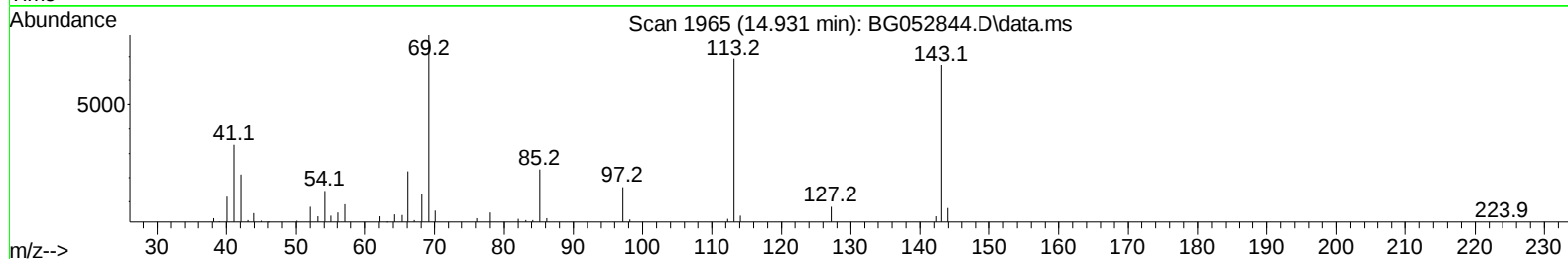
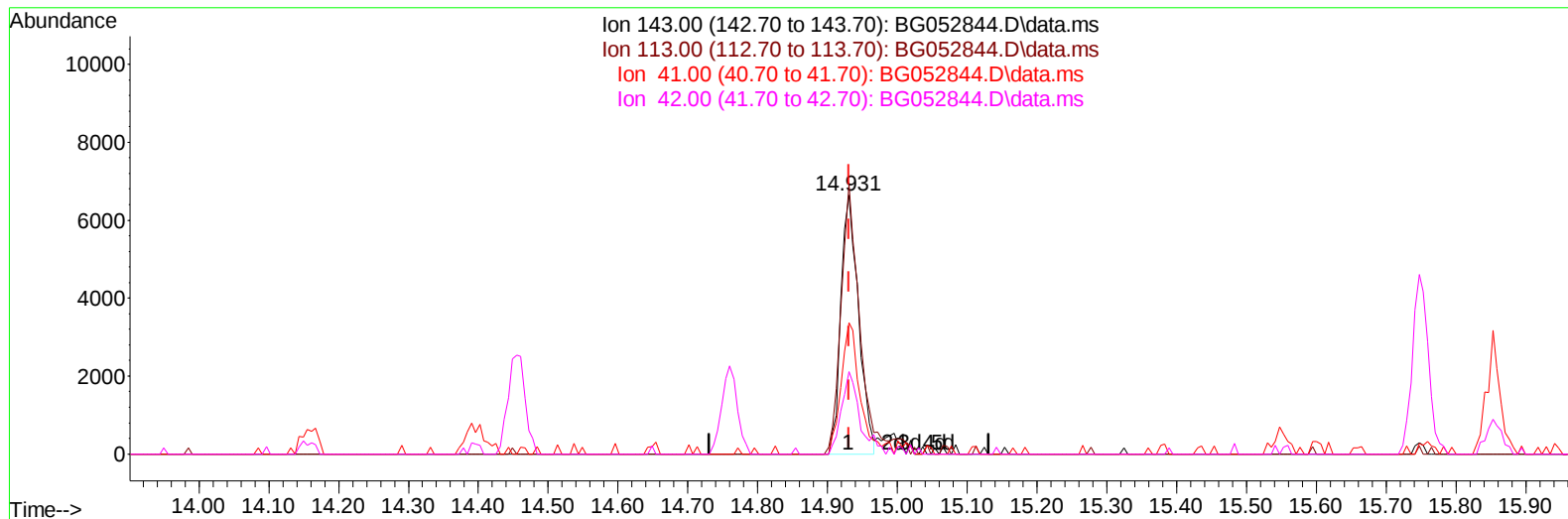
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TIC: BG052844.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.001) 9.25 ng/u1

response 11614

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.19
41.00	47.20	50.66
42.00	29.40	31.86

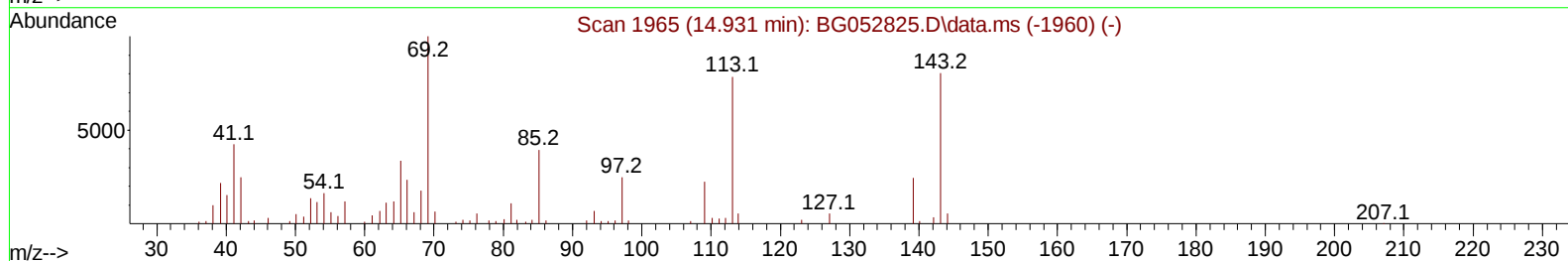
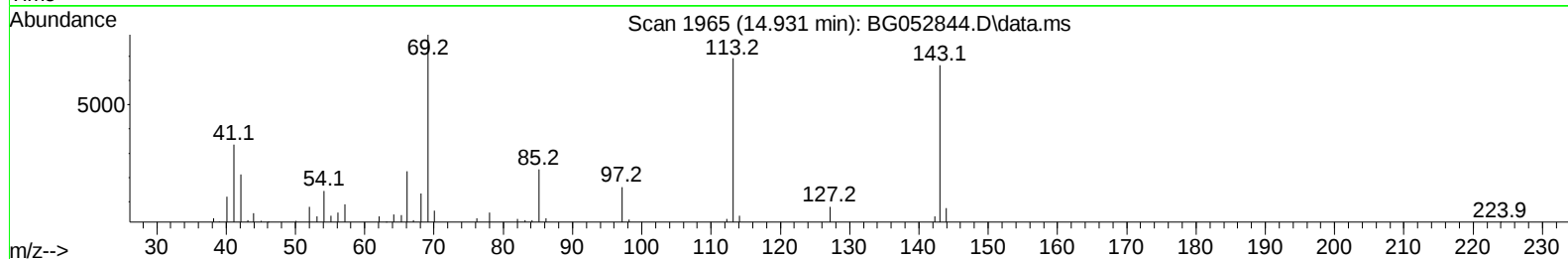
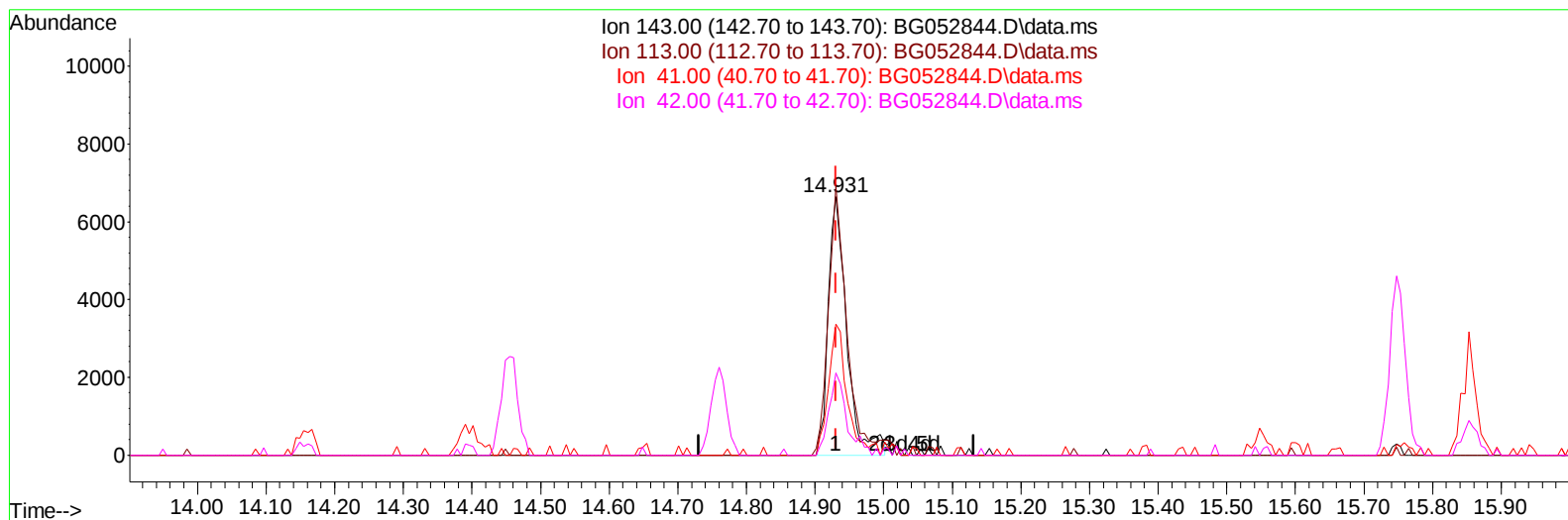
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
 Data File : BG052844.D
 Acq On : 26 Mar 2022 1:09
 Operator : CG/JU
 Sample : N2079-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BG052844.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.001) 9.94 ng/ul m

response 12480

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.19
41.00	47.20	50.66
42.00	29.40	31.86

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	26169	20.000	ng/ul	0.00
20) Naphthalene-d8	10.954	136	117354	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.760	164	97213	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	234405	20.000	ng/ul	0.00
79) Chrysene-d12	21.792	240	261767	20.000	ng/ul	0.00
88) Perylene-d12	25.111	264	262129	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.558	96	2953	3.986	ng/uL	0.00
4) Pyridine-d5	3.975	84	13622m	7.365	ng/ul	0.00
7) Phenol-d5	7.288	99	20545	8.766	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.458	67	53419	37.317	ng/ul	0.00
11) 2-Chlorophenol-d4	7.670	132	45224	28.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.833	113	34517	20.330	ng/ul	0.00
21) Nitrobenzene-d5	9.297	128	32206	38.124	ng/ul	0.00
24) 2-Nitrophenol-d4	10.025	143	36475	40.937	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.566	165	61411	31.681	ng/ul	0.00
31) 4-Chloroaniline-d4	11.077	131	82568	32.229	ng/ul	0.00
46) Dimethylphthalate-d6	14.155	166	281729	37.723	ng/ul	0.00
49) Acenaphthylene-d8	14.455	160	320946	35.365	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	12480m	9.939	ng/ul	0.00
60) Fluorene-d10	15.753	176	248382	37.877	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.853	200	48927	41.016	ng/ul	0.00
73) Anthracene-d10	17.603	188	450591	41.281	ng/ul	0.00
81) Pyrene-d10	19.889	212	563772	38.499	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.882	264	578303	42.697	ng/ul	0.00

Target Compounds Qvalue

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

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