

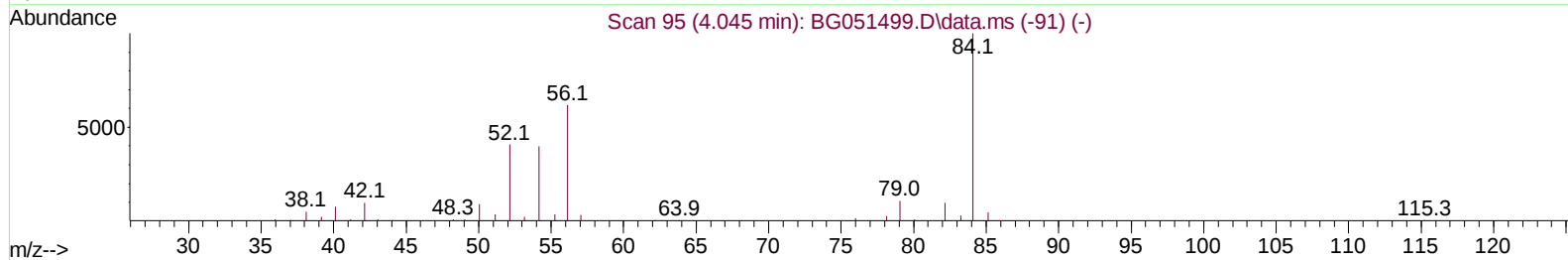
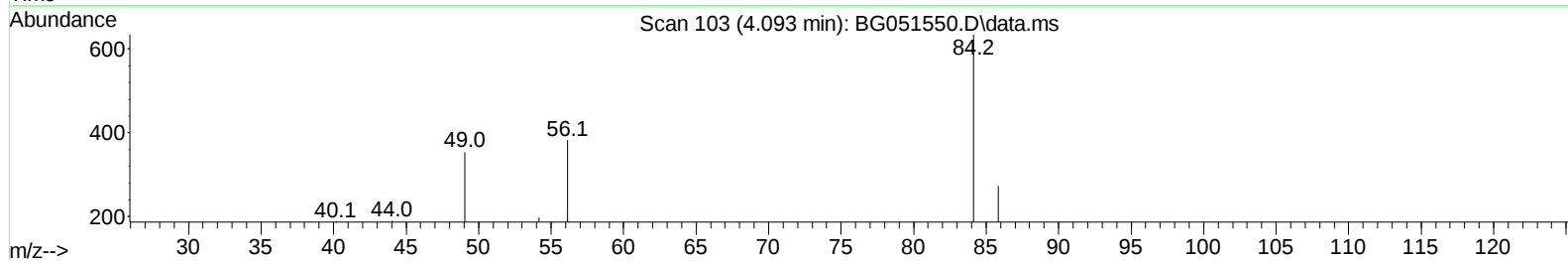
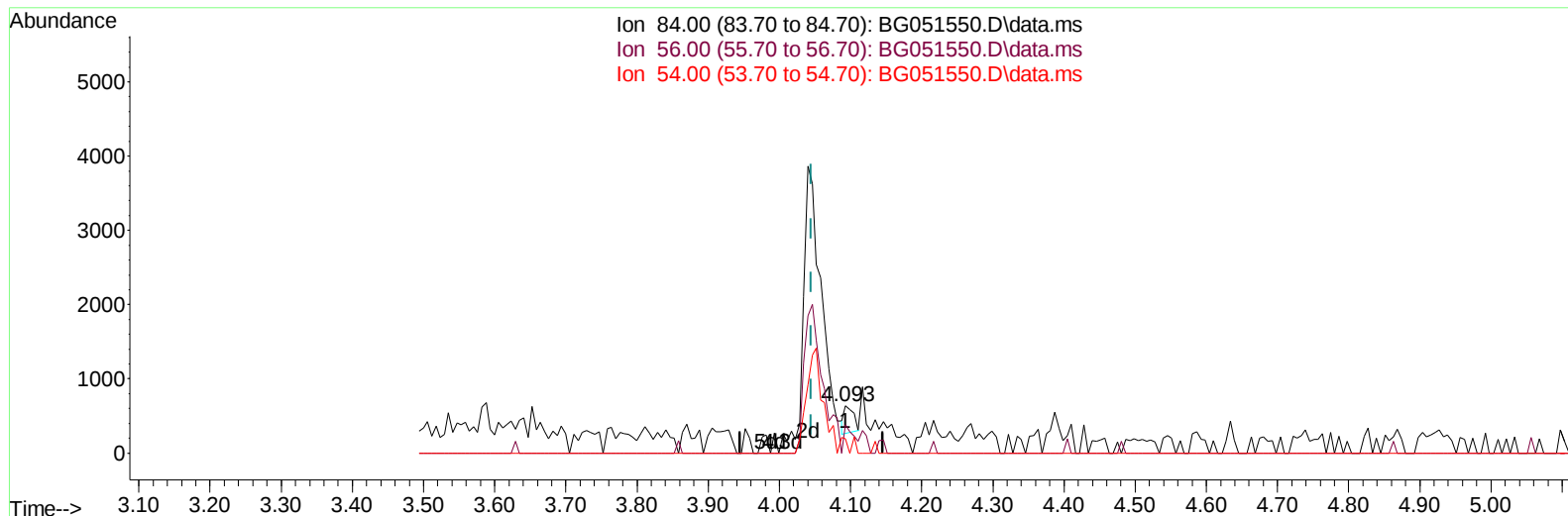
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051550.D
 Acq On : 15 Dec 2021 23:48
 Operator : CG/JU
 Sample : M4995-08
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW5Q1

Manual IntegrationsAPPROVED

Quant Time: Dec 16 02:16:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/16/2021
 Supervised By :Yogesh Patel 12/23/2021



TIC: BG051550.D\data.ms

(4) Pyridine-d5 (S)

4.093min (+ 0.048) 0.15 ng/u1

response 335

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	60.25
54.00	31.50	31.07
0.00	0.00	0.00

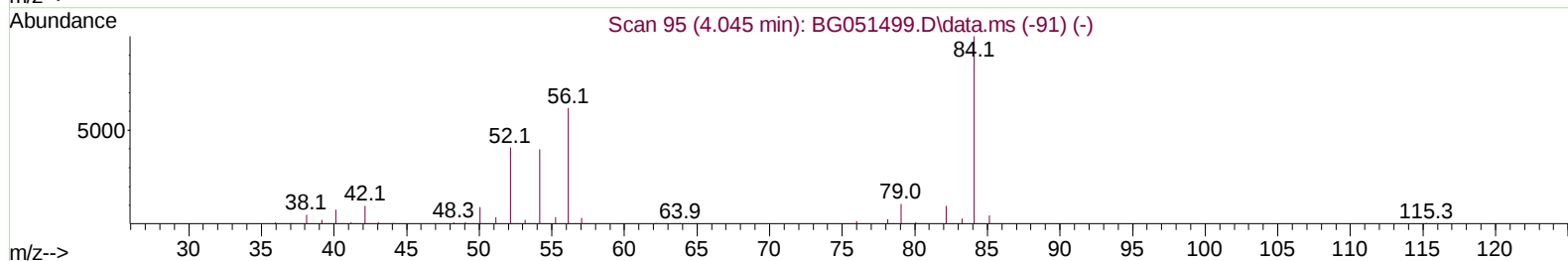
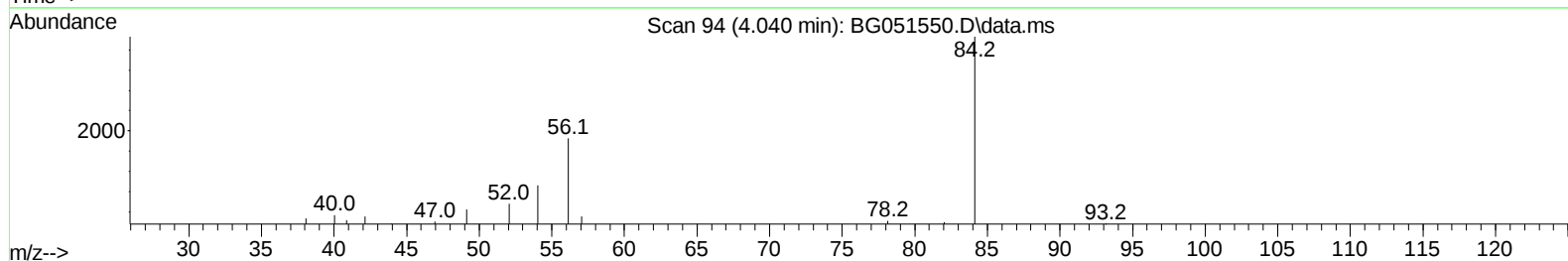
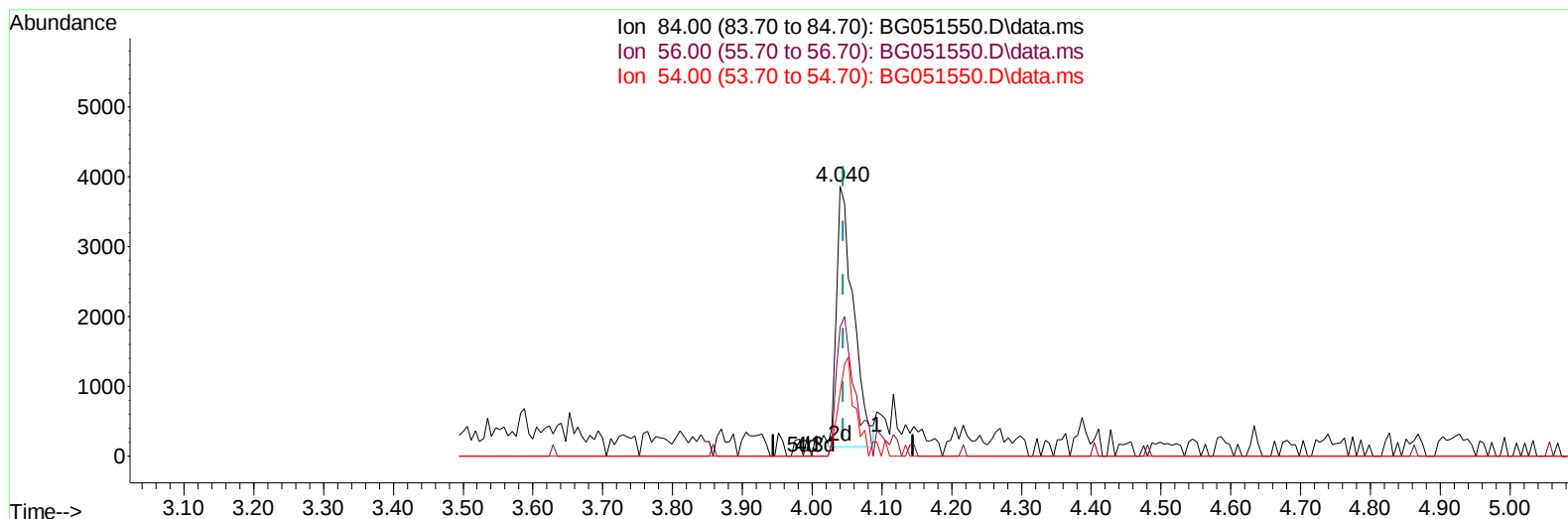
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TIC: BG051550.D\data.ms

(4) Pyridine-d5 (S)

4.040min (-0.005) 2.79 ng/ul m

response 6305

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	47.85#
54.00	31.50	23.80#
0.00	0.00	0.00

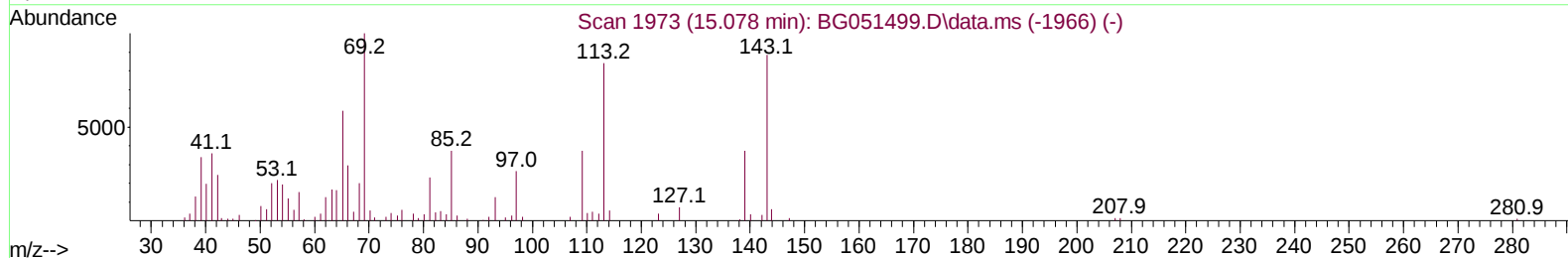
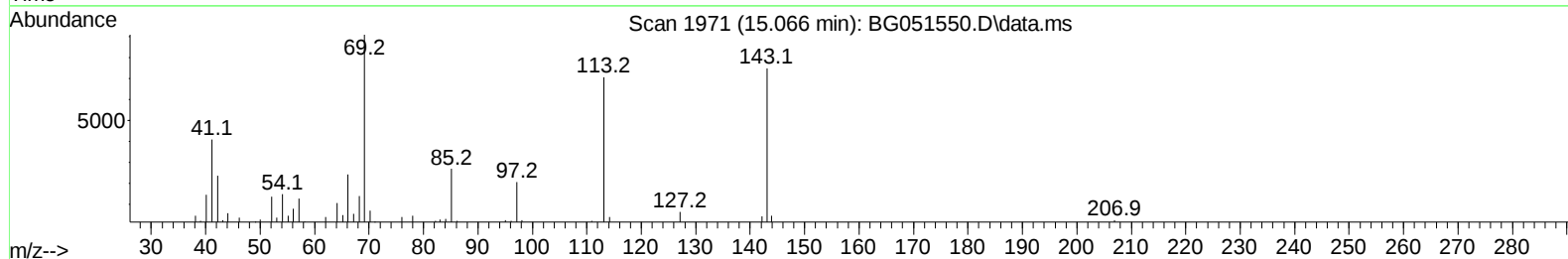
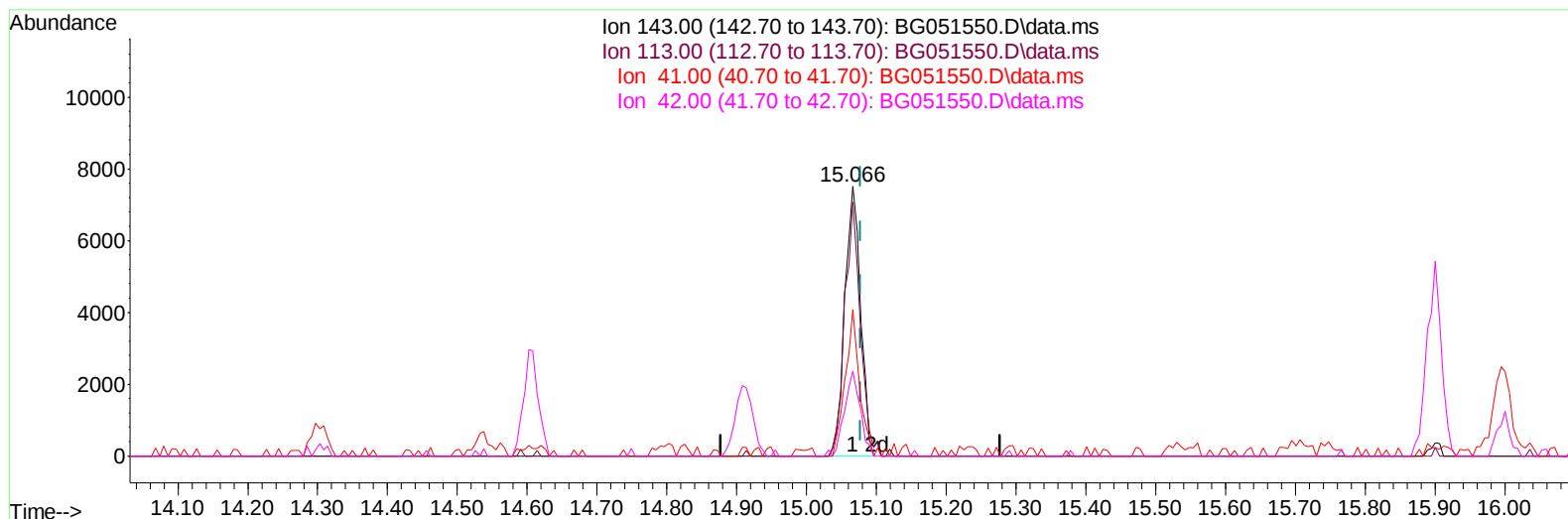
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
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Misc :
ALS Vial : 48 Sample Multiplier: 1

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TIC: BG051550.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (-0.011) 9.99 ng/u1

response 11789

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.27
41.00	44.40	54.45#
42.00	29.70	31.42

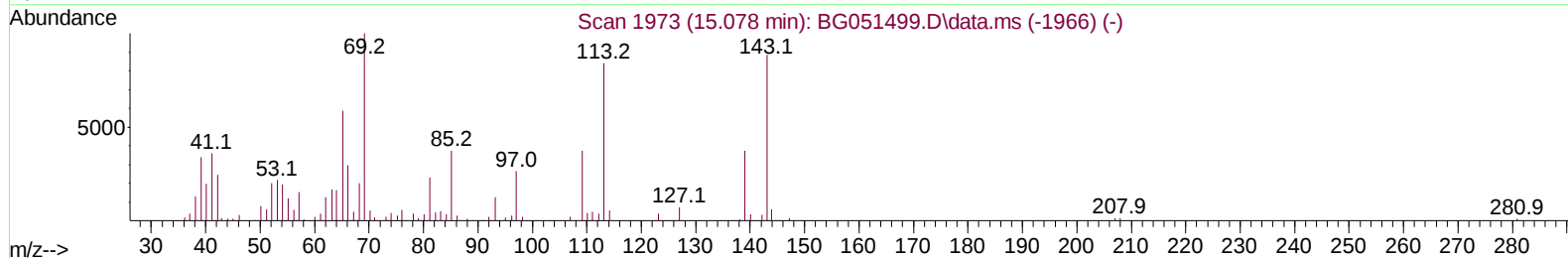
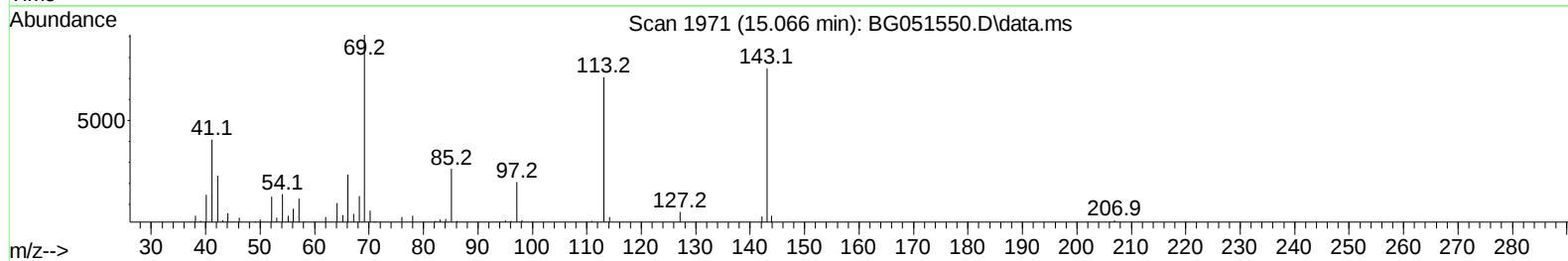
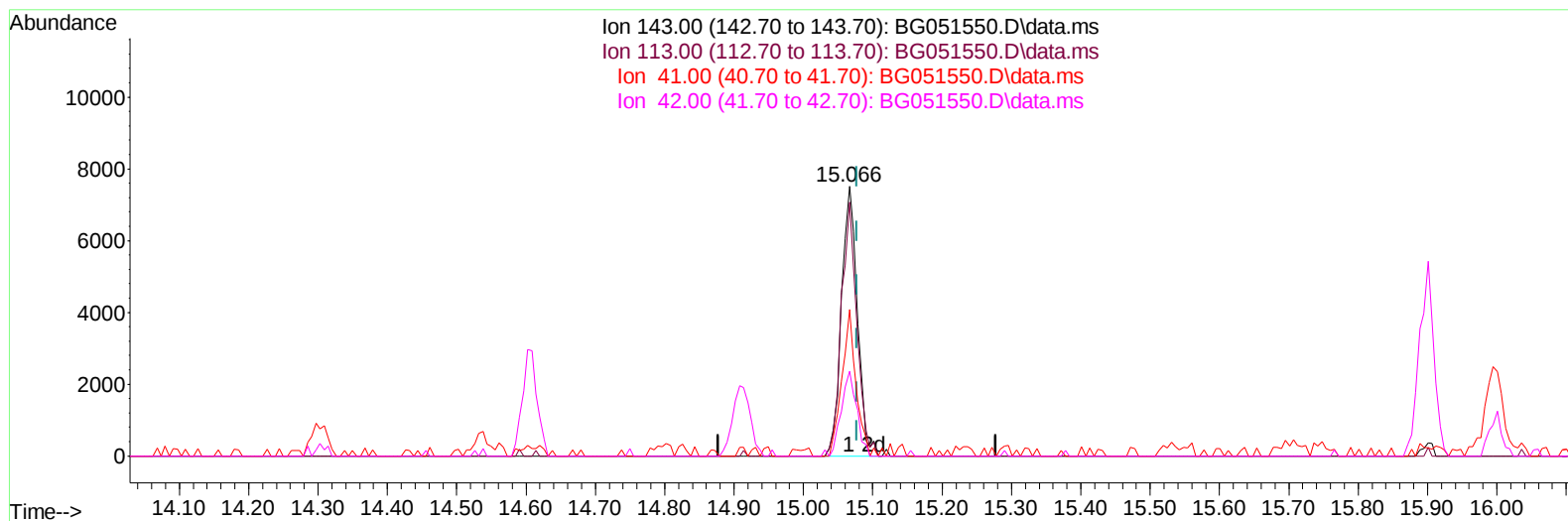
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 Operator : CG/JU
 Sample : M4995-08
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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TIC: BG051550.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (-0.011) 10.12 ng/ul m

response 11938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.27
41.00	44.40	54.45#
42.00	29.70	31.42

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.276	152		31023	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.113	136		124295	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.914	164		90661	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.657	188		226003	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.975	240		233634	20.000	ng/ul	#-0.01
88) Perylene-d12	25.470	264		234306	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.606	96		4080	5.462	ng/uL	0.00
4) Pyridine-d5	4.040	84		6305m	2.794	ng/ul	0.00
7) Phenol-d5	7.406	99		22789	8.227	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.582	67		57341	34.822	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.806	132		52719	27.365	ng/ul	0.00
15) 4-Methylphenol-d8	8.969	113		42485	19.842	ng/ul	-0.01
21) Nitrobenzene-d5	9.445	128		34343	37.687	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.179	143		37749	37.894	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.719	165		71883	33.275	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.236	131		65979	23.197	ng/ul	0.00
46) Dimethylphthalate-d6	14.303	166		305844	42.148	ng/ul	-0.01
49) Acenaphthylene-d8	14.608	160		337898	37.493	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143		11938m	10.117	ng/ul	-0.01
60) Fluorene-d10	15.901	176		267110	41.086	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.000	200		53530	45.461	ng/ul	-0.01
73) Anthracene-d10	17.757	188		482844	44.787	ng/ul	0.00
81) Pyrene-d10	20.030	212		599326	42.622	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.235	264		568210	46.090	ng/ul	-0.02
Target Compounds							
						Qvalue	
2) 1,4-Dioxane	3.641	88		7132	9.148	ng/ul#	91

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

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