

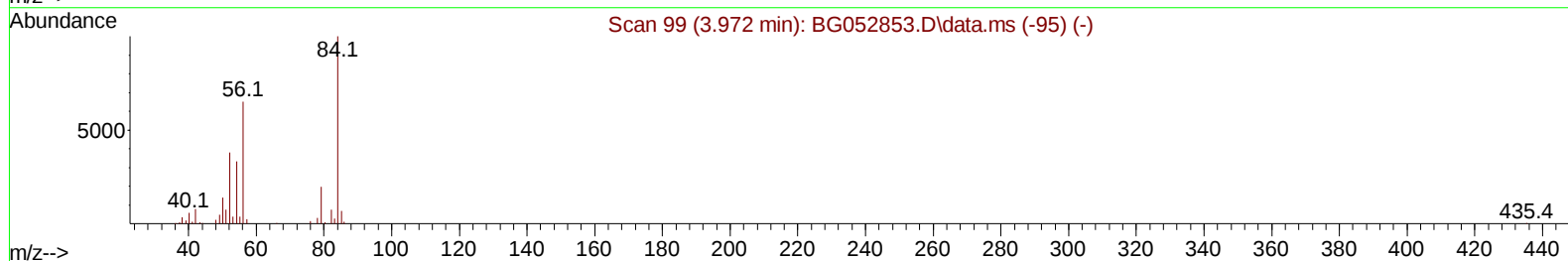
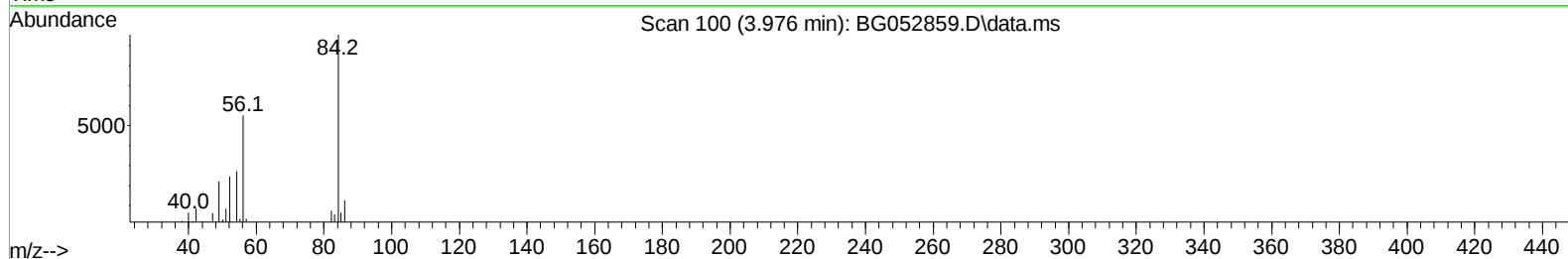
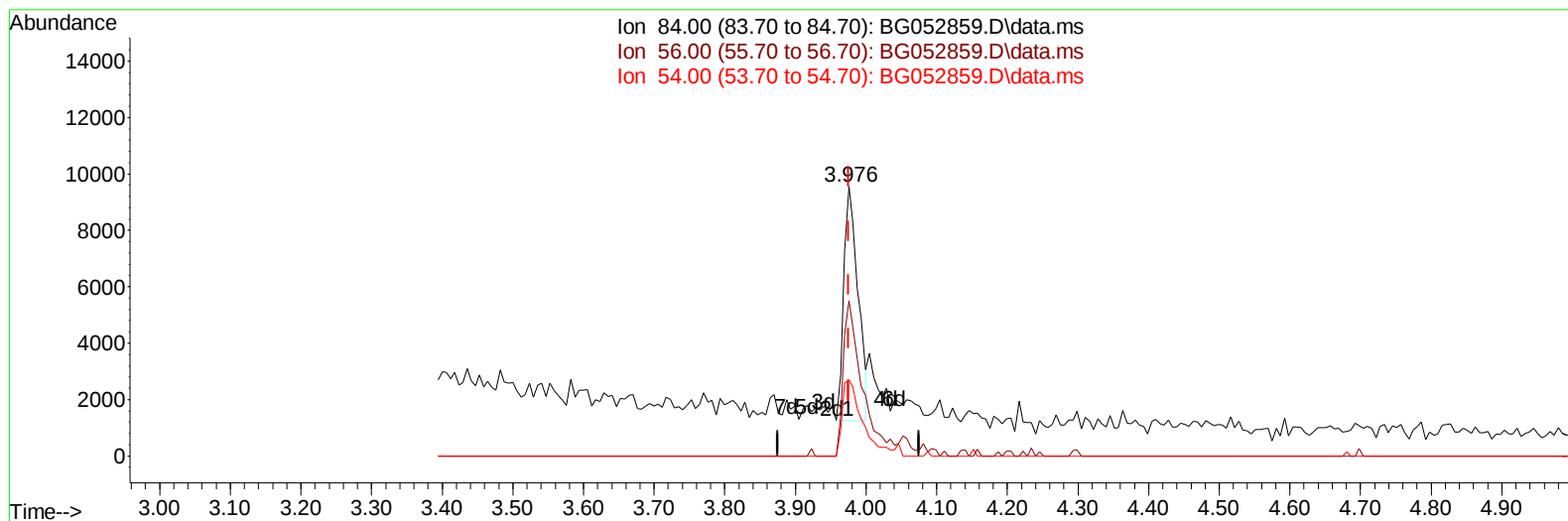
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
 Data File : BG052859.D
 Acq On : 26 Mar 2022 11:31
 Operator : CG/JU
 Sample : N2068-02
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YAZ46

Manual IntegrationsAPPROVED

Quant Time: Mar 28 00:58:13 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: BG052859.D\data.ms

(4) Pyridine-d5 (S)

3.976min (+ 0.000) 6.31 ng/u1

response 11658

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	57.71
54.00	28.70	28.27
0.00	0.00	0.00

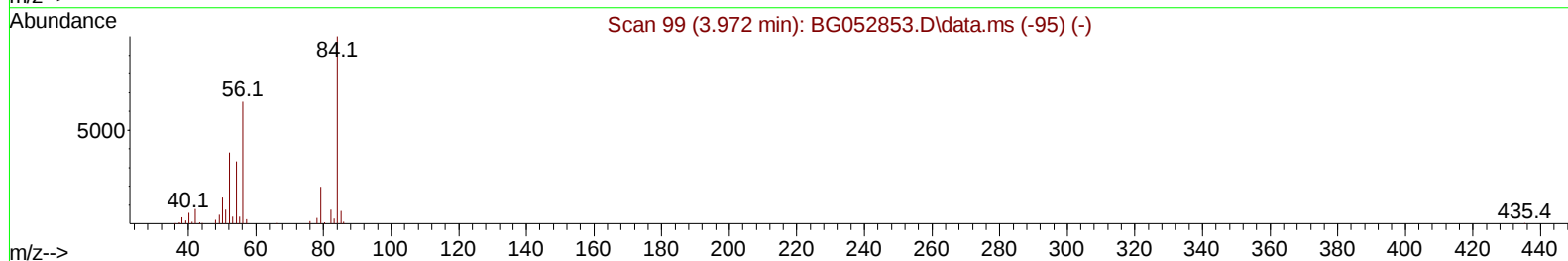
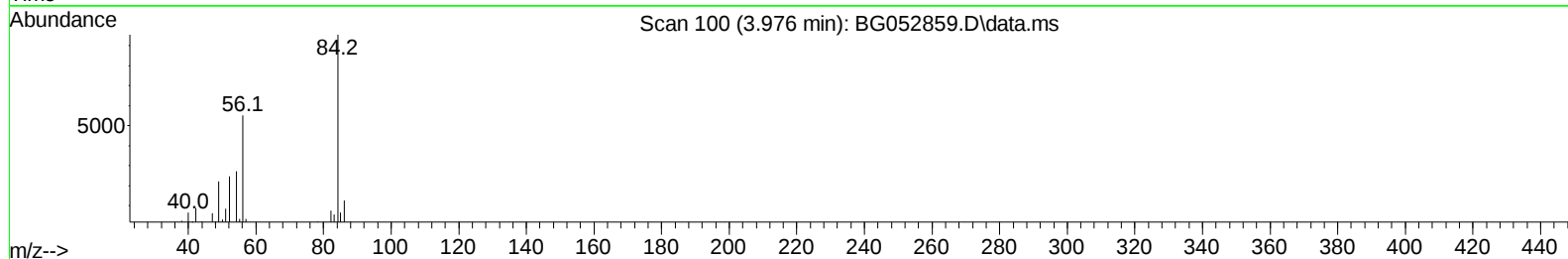
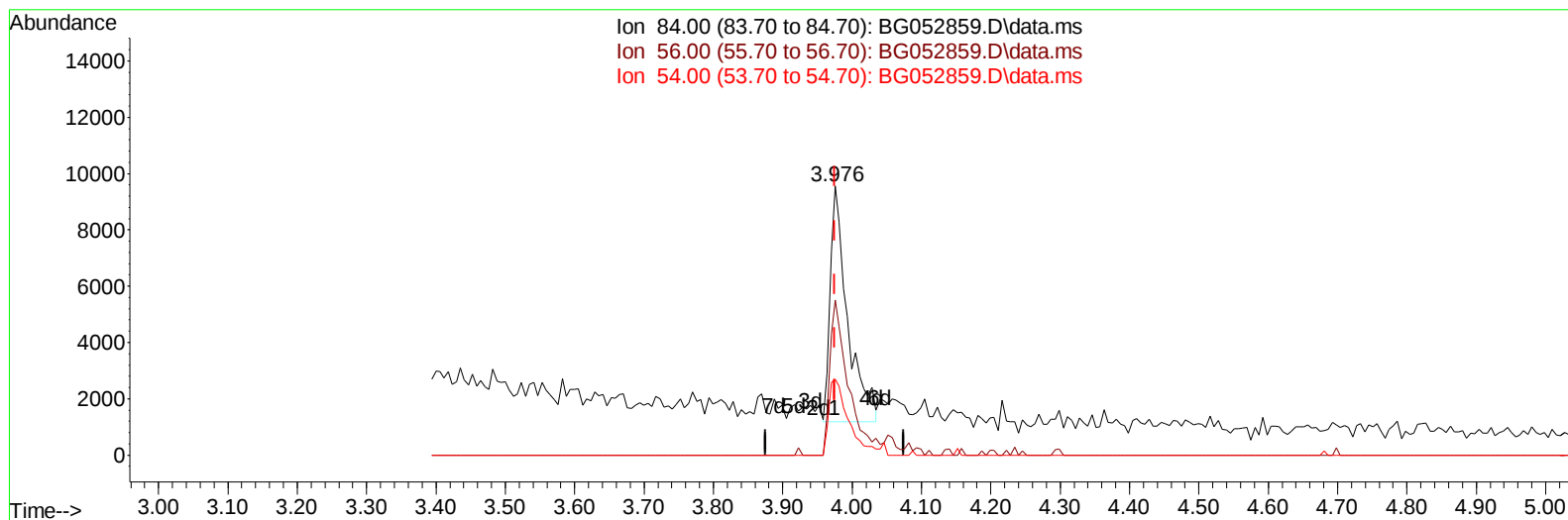
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(4) Pyridine-d5 (S)

3.976min (+ 0.000) 7.87 ng/u1 m

response 14546

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	57.71
54.00	28.70	28.27
0.00	0.00	0.00

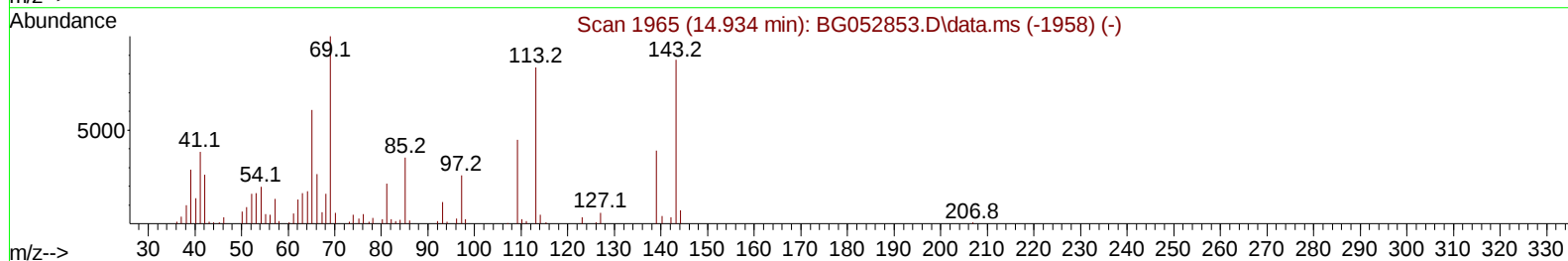
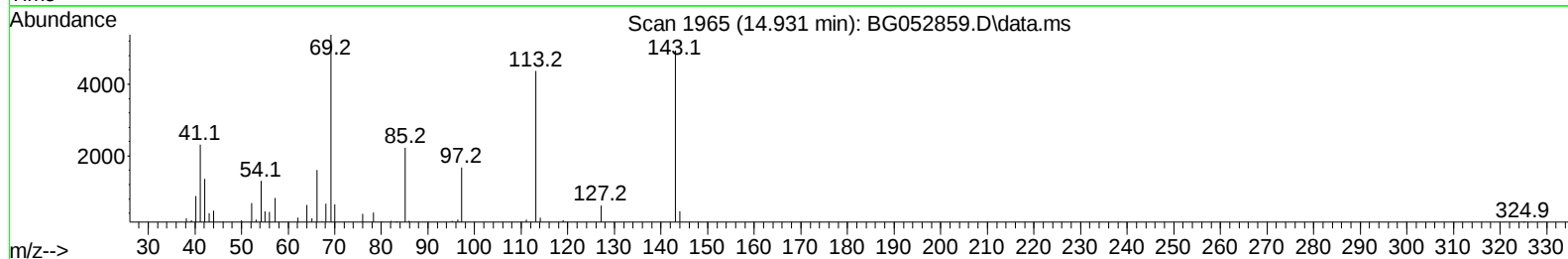
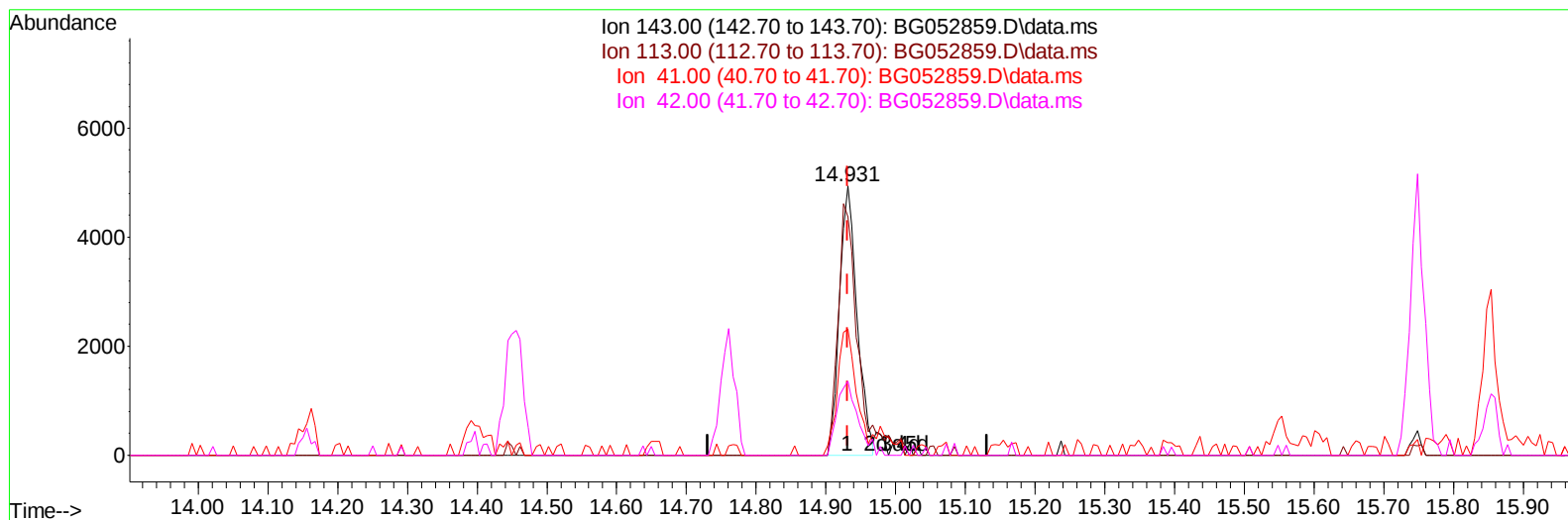
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TIC: BG052859.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (+ 0.000) 6.82 ng/u1

response 8624

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	88.47
41.00	47.20	46.79
42.00	29.40	27.53

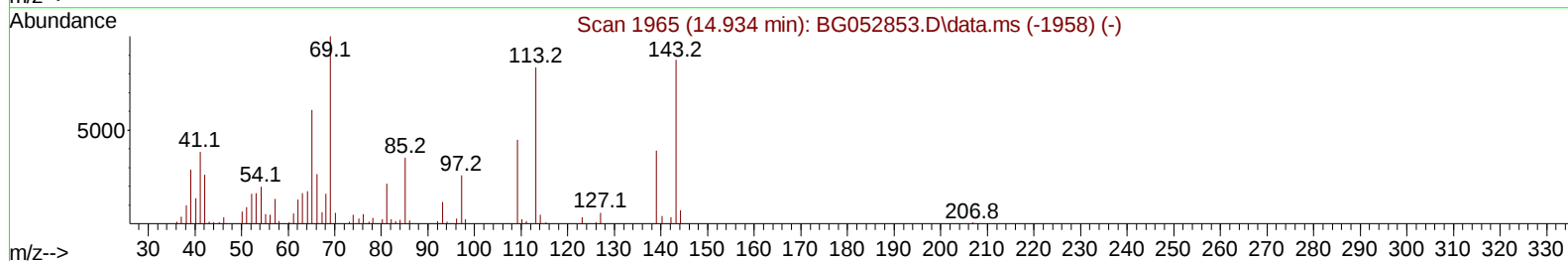
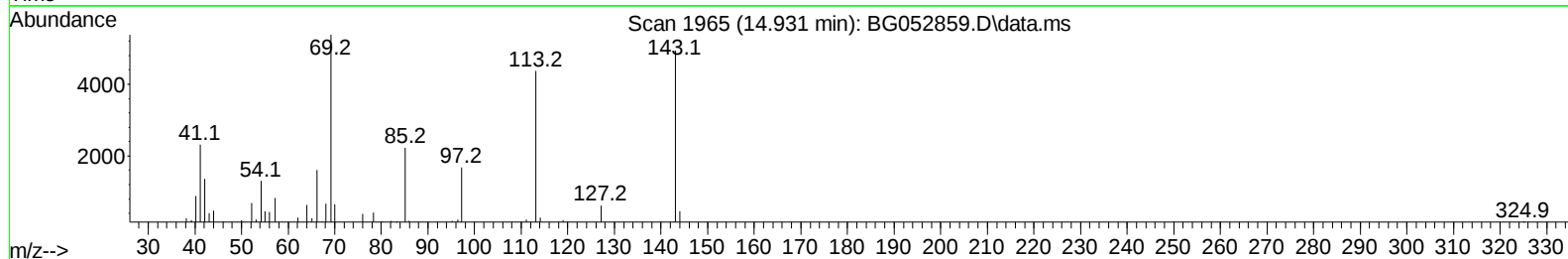
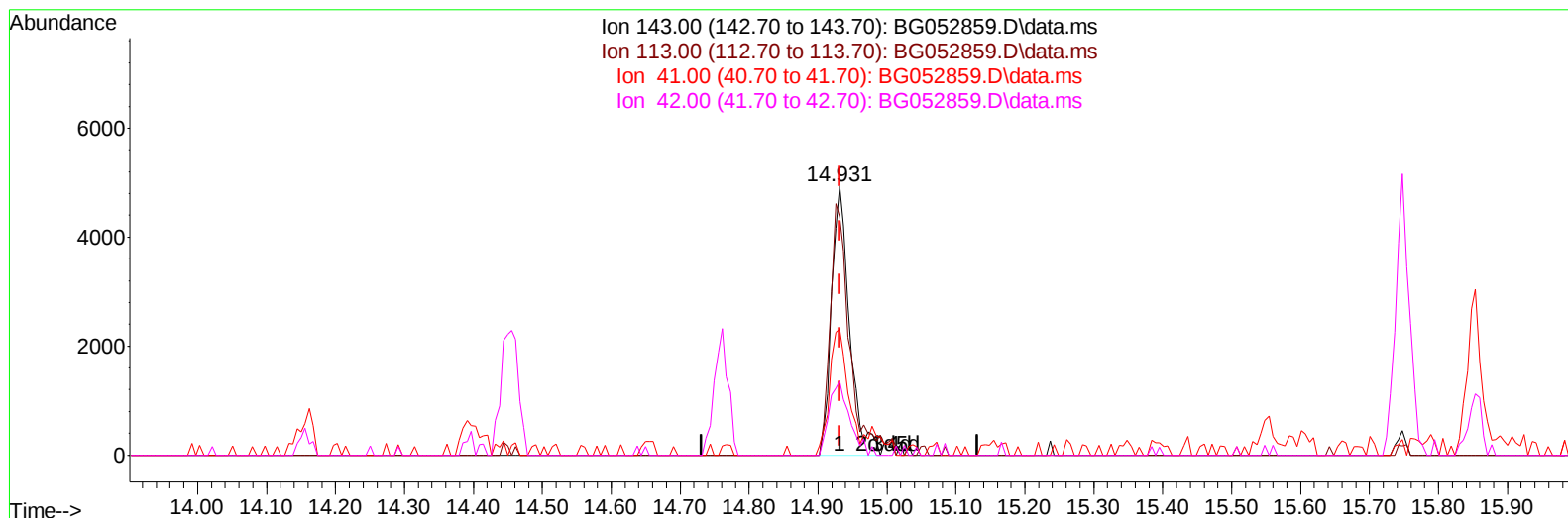
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Operator : CG/JU
Sample : N2068-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG052859.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (+ 0.000) 7.11 ng/ul m

response 8997

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	88.47
41.00	47.20	46.79
42.00	29.40	27.53

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.141	152	26161	20.000	ng/ul	0.00
20) Naphthalene-d8	10.954	136	114858	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.761	164	97908	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	239665	20.000	ng/ul	0.00
79) Chrysene-d12	21.787	240	267165	20.000	ng/ul	0.00
88) Perylene-d12	25.106	264	269706	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.553	96	2672	3.608	ng/uL	0.00
4) Pyridine-d5	3.976	84	14546m	7.867	ng/ul	0.00
7) Phenol-d5	7.283	99	16670	7.114	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.459	67	52195	36.473	ng/ul	0.00
11) 2-Chlorophenol-d4	7.671	132	41362	25.998	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	31961	18.830	ng/ul	0.00
21) Nitrobenzene-d5	9.298	128	32557	39.377	ng/ul	0.00
24) 2-Nitrophenol-d4	10.026	143	34145	39.154	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.567	165	62461	32.923	ng/ul	0.00
31) 4-Chloroaniline-d4	11.078	131	66695	26.599	ng/ul	0.00
46) Dimethylphthalate-d6	14.156	166	271320	36.072	ng/ul	0.00
49) Acenaphthylene-d8	14.456	160	310555	33.978	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	8997m	7.114	ng/ul	0.00
60) Fluorene-d10	15.748	176	236354	35.787	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.854	200	46561	38.176	ng/ul	0.00
73) Anthracene-d10	17.604	188	418585	37.507	ng/ul	0.00
81) Pyrene-d10	19.883	212	534121	35.737	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.883	264	552849	39.671	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.312	94	2821	1.249	ng/ul#	86
78) Di-n-butylphthalate	18.456	149	20287	1.535	ng/ul#	96

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

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