

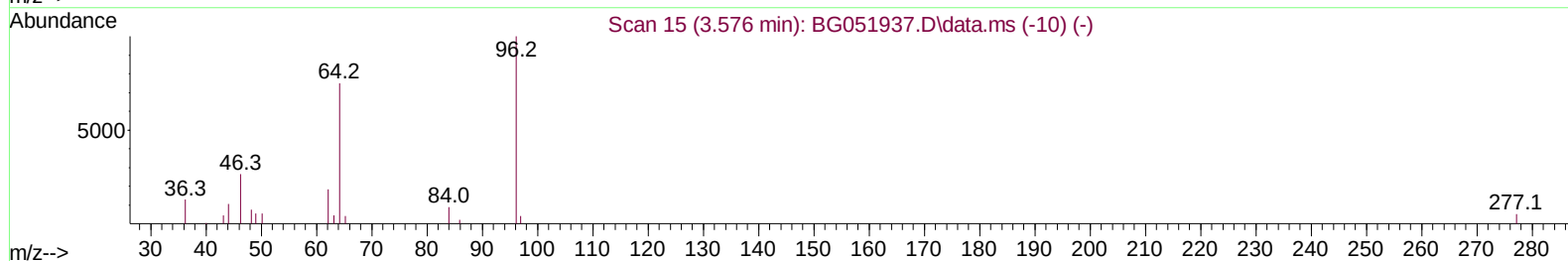
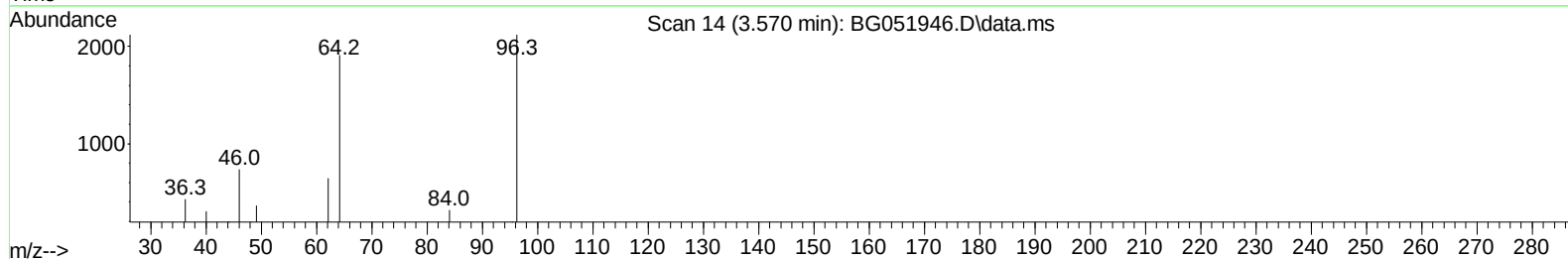
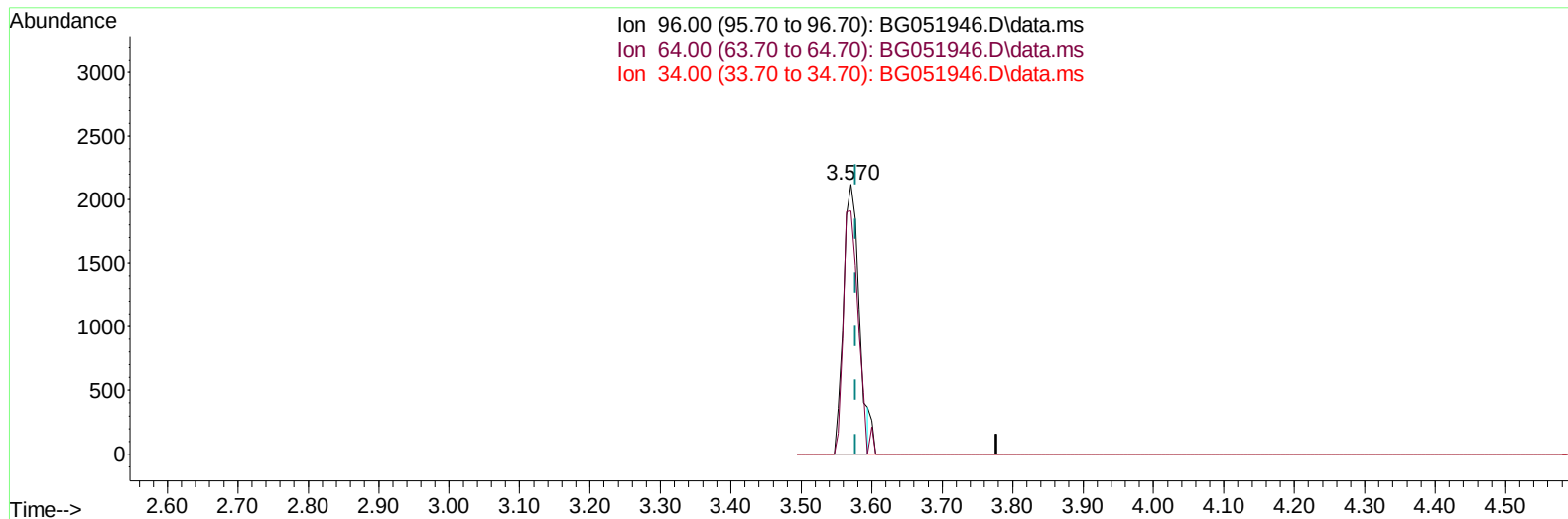
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011222\
 Data File : BG051946.D
 Acq On : 12 Jan 2022 15:20
 Operator : CG/JU
 Sample : N1100-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLN7

Manual IntegrationsAPPROVED

Quant Time: Jan 12 23:53:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/13/2022
 Supervised By :mohammad ahmed 01/18/2022



TIC: BG051946.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.570min (-0.007) 5.07 ng/uL

response 3196

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	90.28
34.00	0.00	0.00
0.00	0.00	0.00

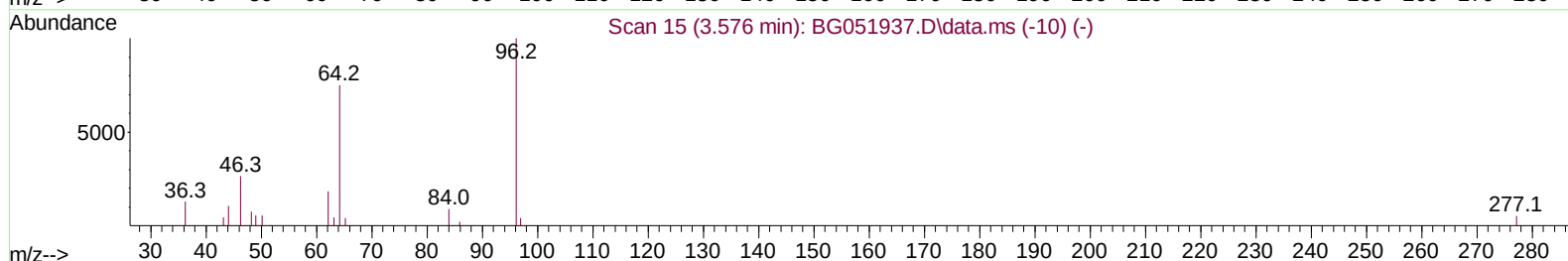
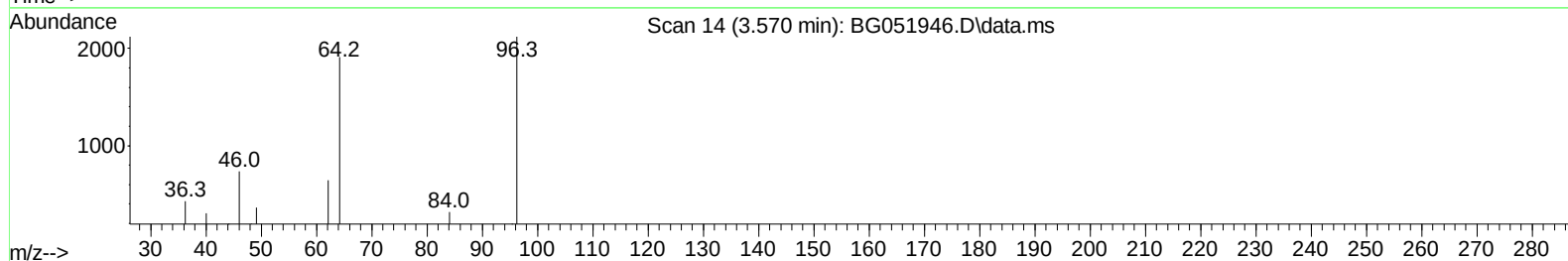
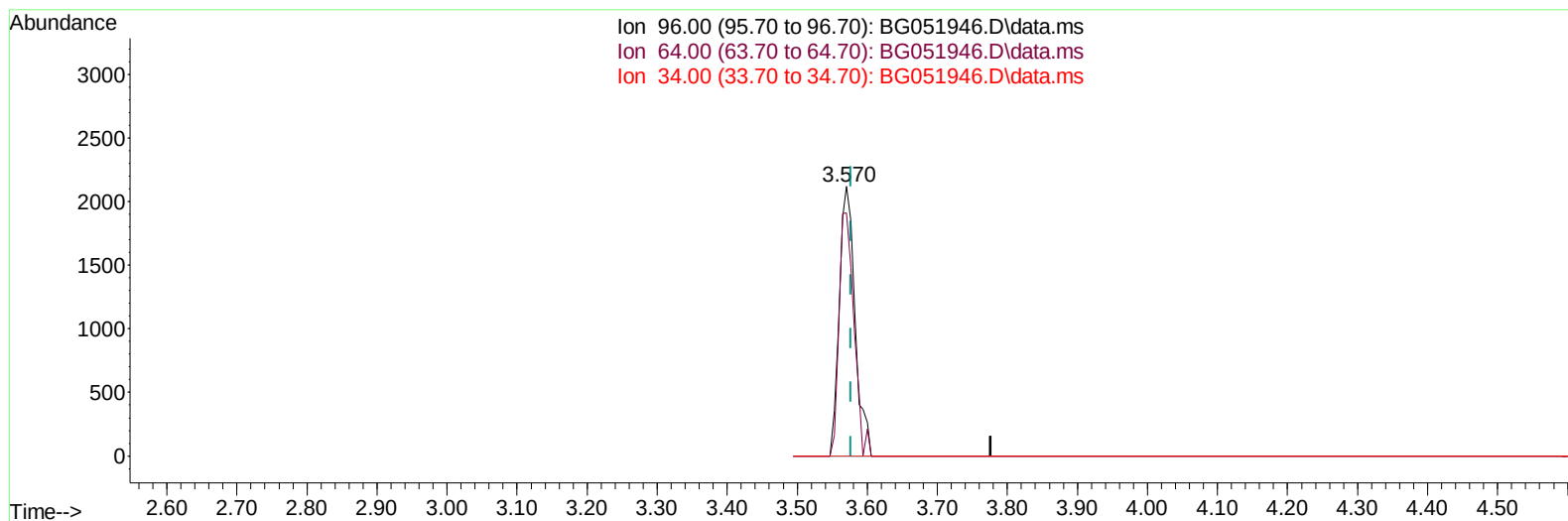
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TIC: BG051946.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.570min (-0.007) 5.21 ng/uL m

response 3289

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	90.28
34.00	0.00	0.00
0.00	0.00	0.00

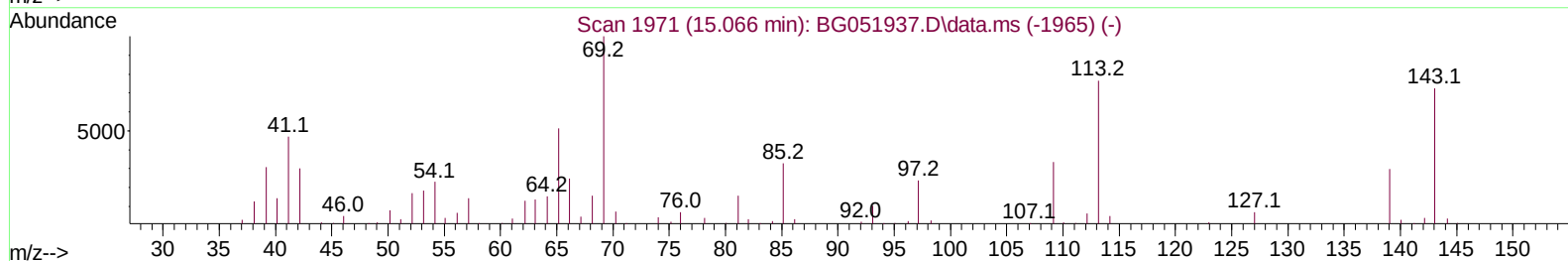
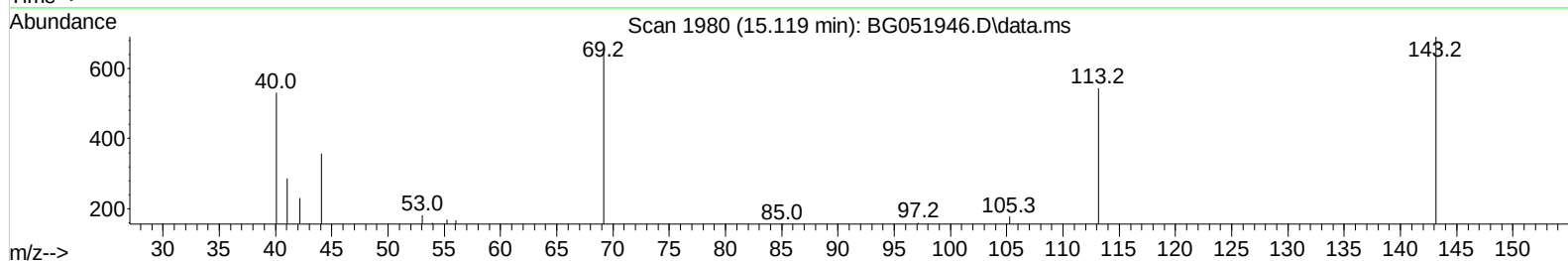
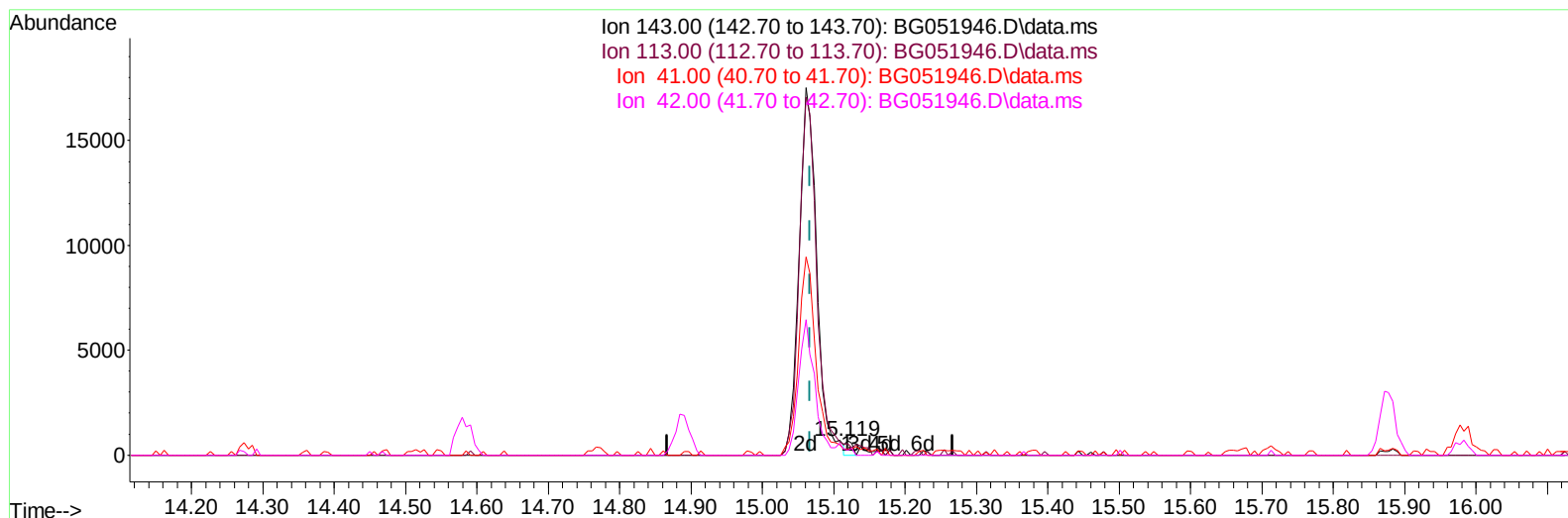
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TIC: BG051946.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.119min (+ 0.052) 0.45 ng/u1

response 440

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	78.87
41.00	44.40	41.53
42.00	29.70	33.43

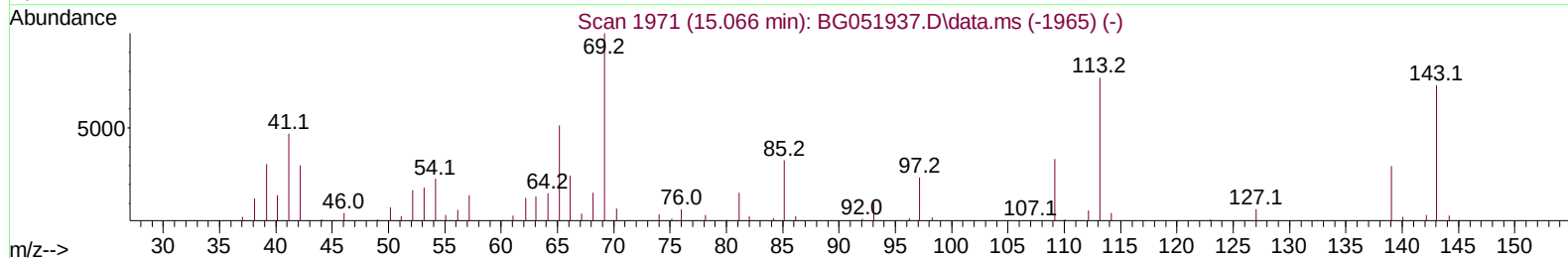
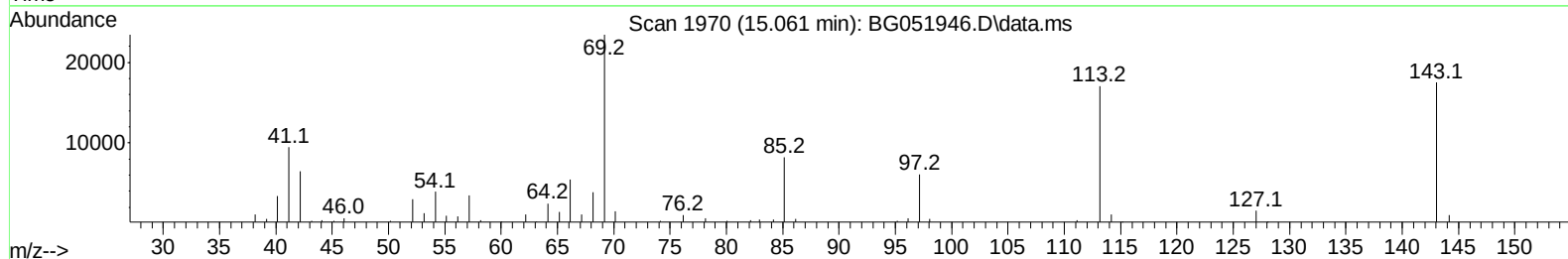
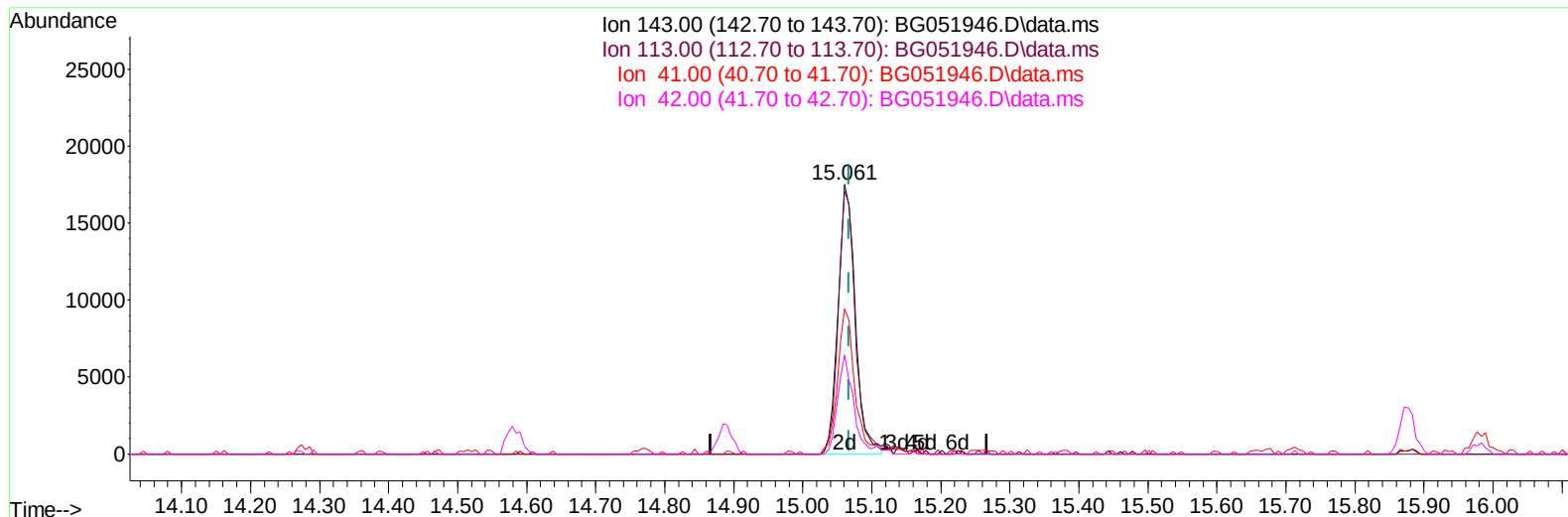
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Instrument :
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TIC: BG051946.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.061min (-0.006) 31.01 ng/ul m

response 30121

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	97.44#
41.00	44.40	53.88#
42.00	29.70	36.79#

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 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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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	8.246	152	22545	20.000	ng/ul	0.00
20) Naphthalene-d8	11.084	136	91252	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.890	164	70282	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.640	188	196303	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.957	240	200327	20.000	ng/ul	# 0.00
88) Perylene-d12	25.441	264	199303	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.570	96	3289m	5.213	ng/uL	0.00
4) Pyridine-d5	3.999	84	39867	20.327	ng/ul	0.00
7) Phenol-d5	7.383	99	61640	27.935	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	38371	27.394	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.771	132	43681	29.900	ng/ul	0.00
15) 4-Methylphenol-d8	8.945	113	45483	26.516	ng/ul	0.00
21) Nitrobenzene-d5	9.415	128	21310	29.554	ng/ul	0.00
24) 2-Nitrophenol-d4	10.150	143	25251	31.664	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.696	165	46657	29.932	ng/ul	0.00
31) 4-Chloroaniline-d4	11.207	131	63396	29.205	ng/ul	0.00
46) Dimethylphthalate-d6	14.279	166	167931	28.907	ng/ul	0.00
49) Acenaphthylene-d8	14.585	160	201748	28.750	ng/ul	0.00
54) 4-Nitrophenol-d4	15.061	143	30121m	31.012	ng/ul	0.00
60) Fluorene-d10	15.877	176	150025	29.620	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.983	200	26788	21.632	ng/ul	0.00
73) Anthracene-d10	17.739	188	277157	29.763	ng/ul	0.00
81) Pyrene-d10	20.013	212	355823	30.863	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.200	264	303868	28.961	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	11.131	128	9315	1.884	ng/ul#	87
36) 2-Methylnaphthalene	12.729	142	20930	5.942	ng/ul	89
37) 1-Methylnaphthalene	12.946	142	7266	2.010	ng/ul#	95

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

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