

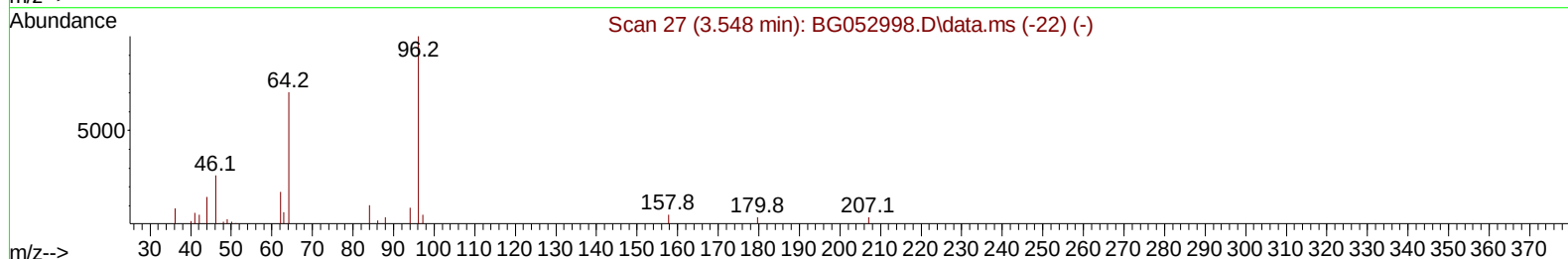
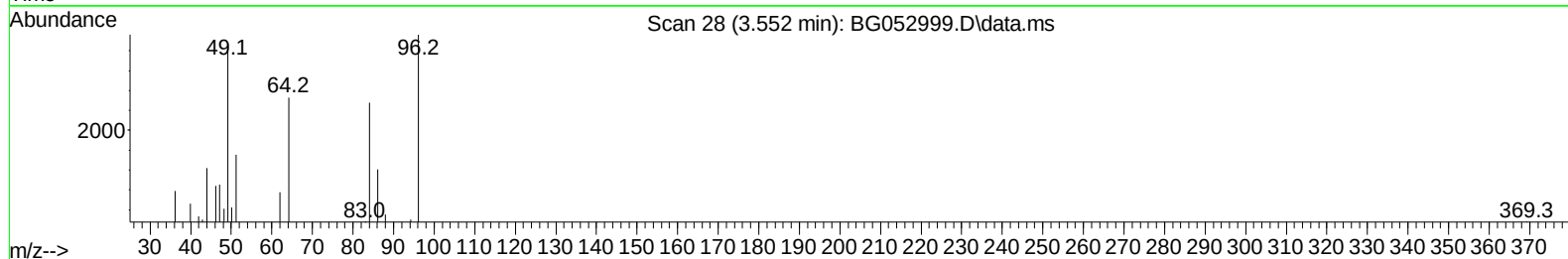
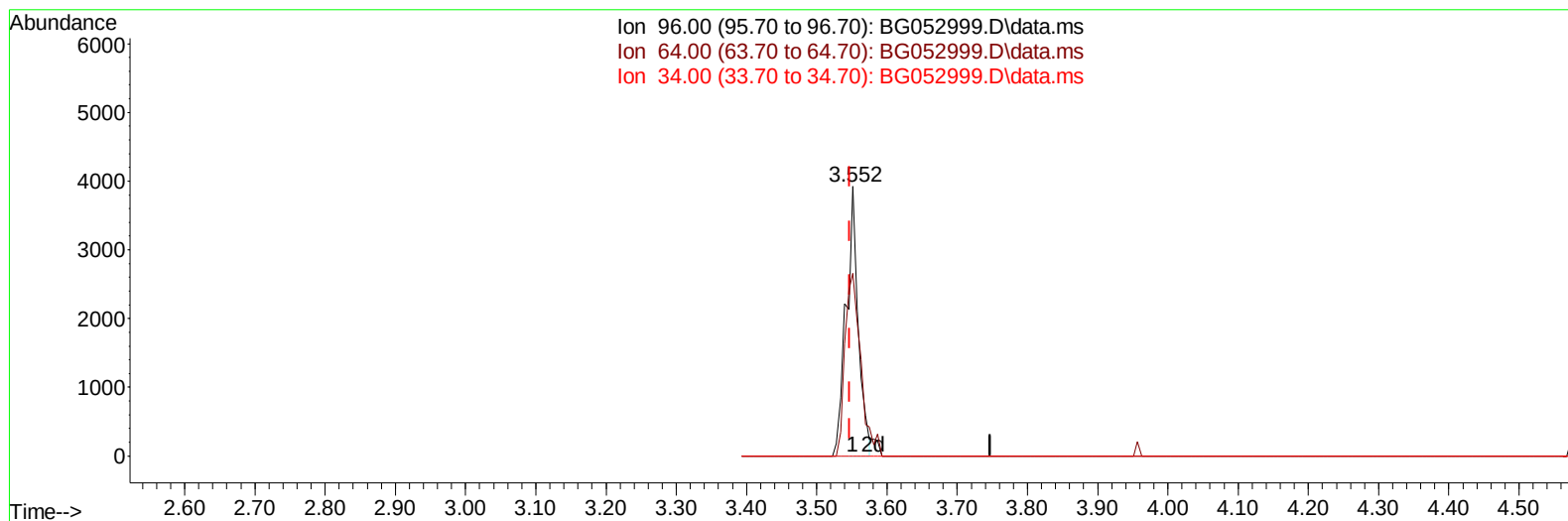
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
 Data File : BG052999.D
 Acq On : 1 Apr 2022 16:16
 Operator : CG/JU
 Sample : PB143815BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK815

Manual IntegrationsAPPROVED

Quant Time: Apr 02 00:37:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 31 00:50:49 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/04/2022
 Supervised By :Yogesh Patel 04/06/2022



TIC: BG052999.D\data.ms

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

3.552min (+ 0.005) 5.43 ng/uL

response 4712

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	67.82
34.00	0.00	0.00
0.00	0.00	0.00

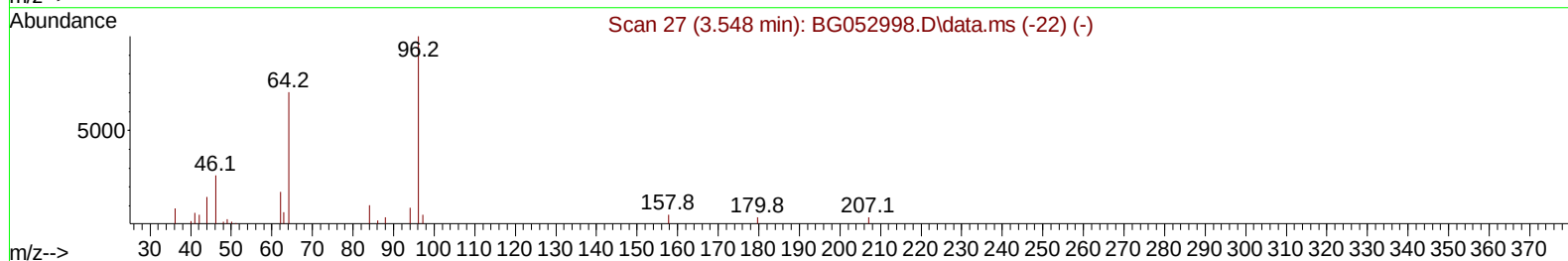
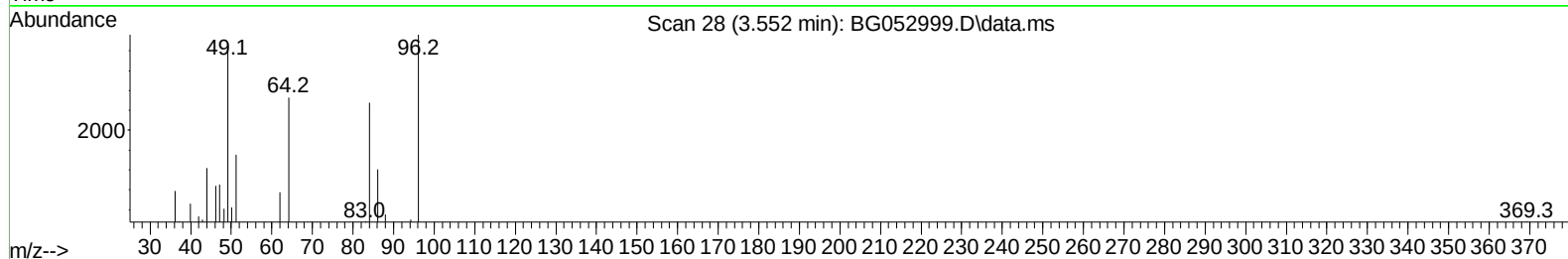
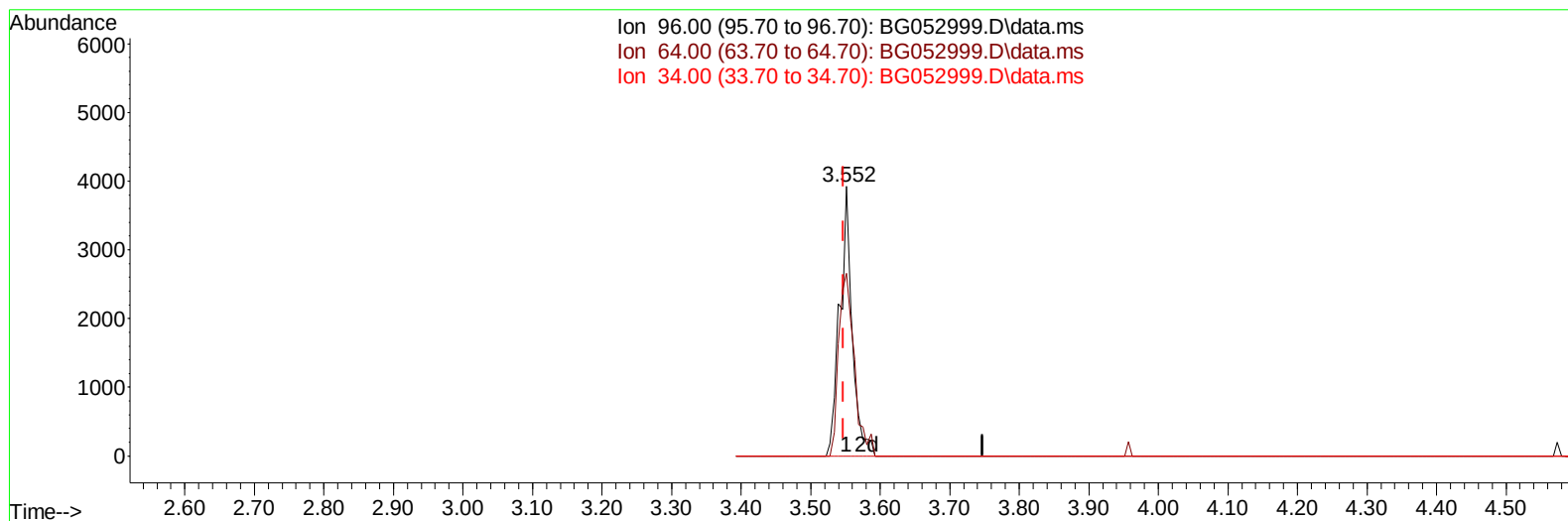
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TIC: BG052999.D\data.ms

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

3.552min (+ 0.005) 5.62 ng/uL m

response 4875

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	67.82
34.00	0.00	0.00
0.00	0.00	0.00

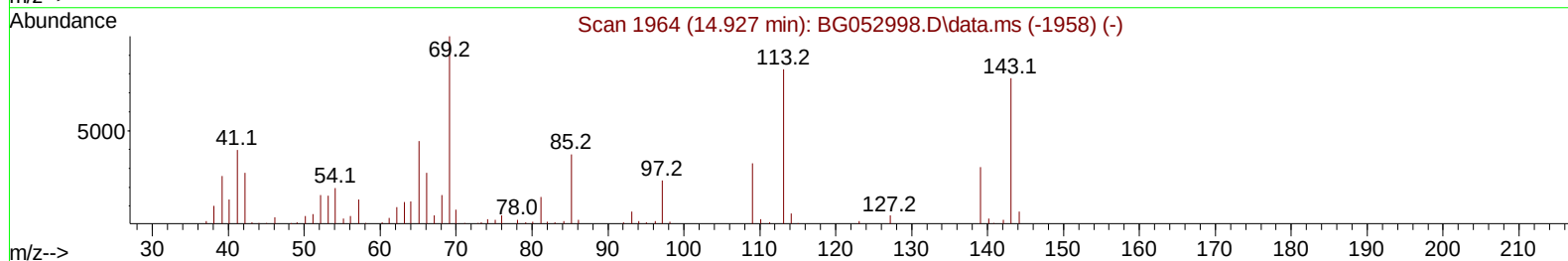
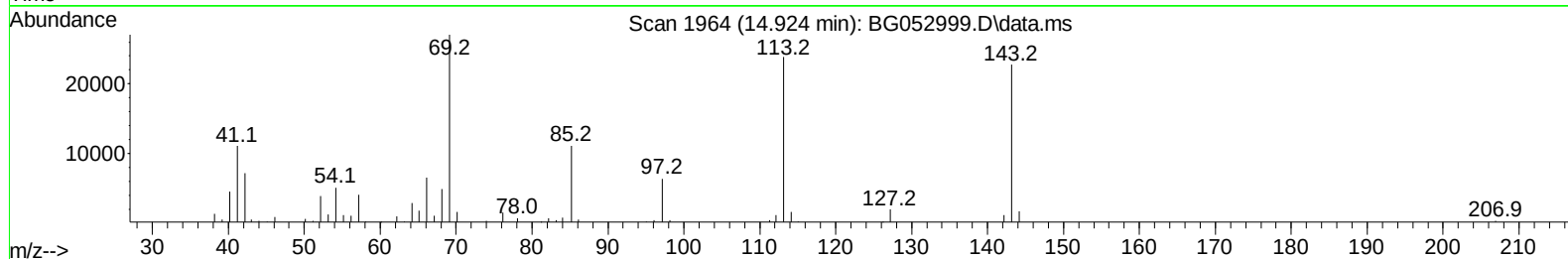
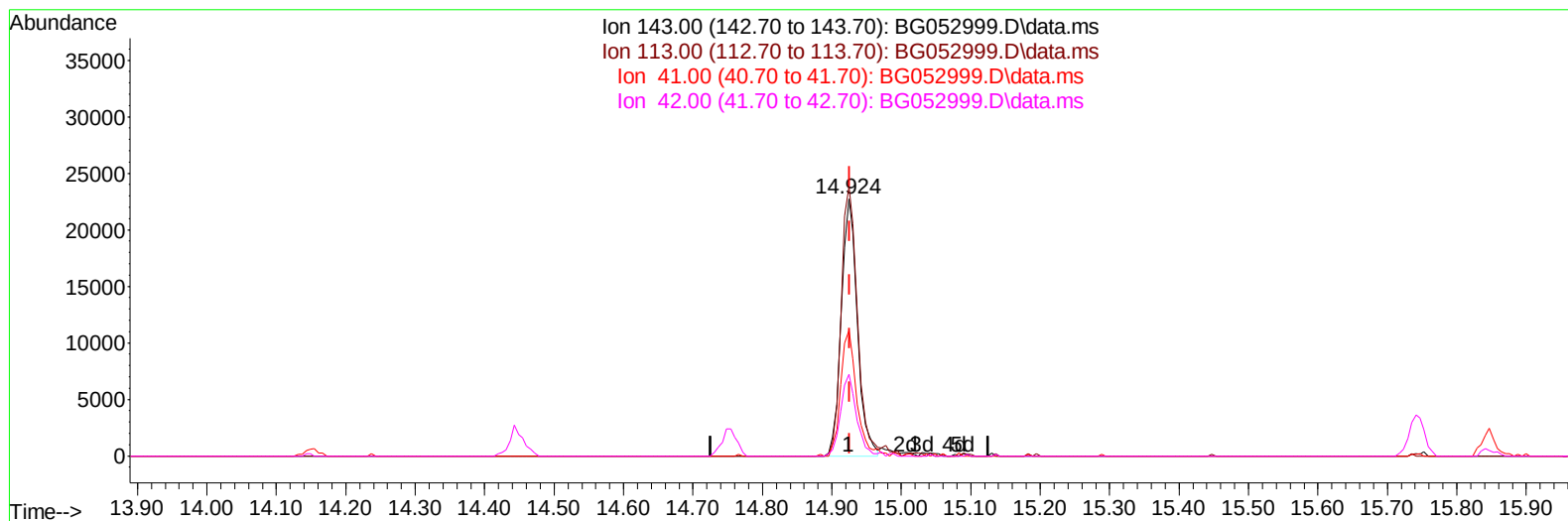
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 Sample : PB143815BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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TIC: BG052999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.924min (-0.001) 28.53 ng/ul

response 36349

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.82
41.00	47.20	48.82
42.00	29.40	31.75

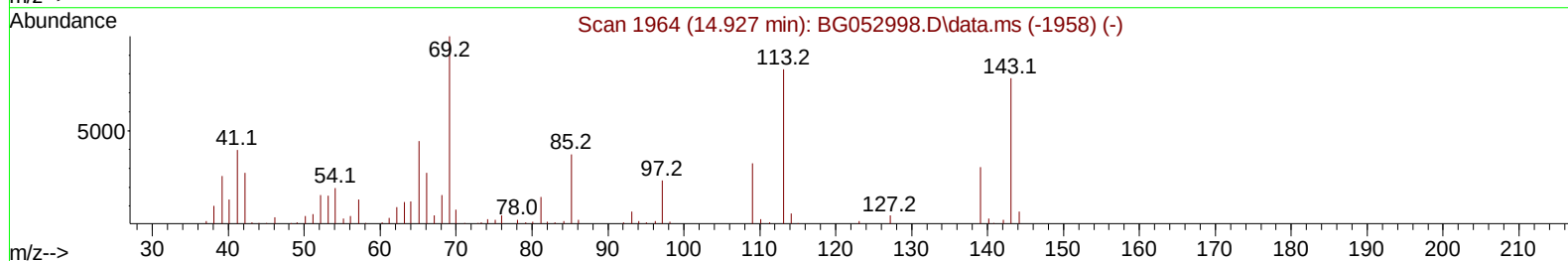
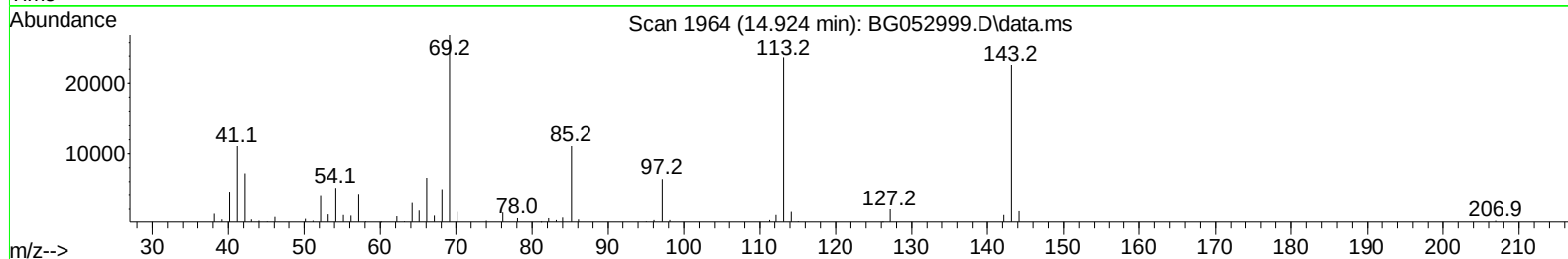
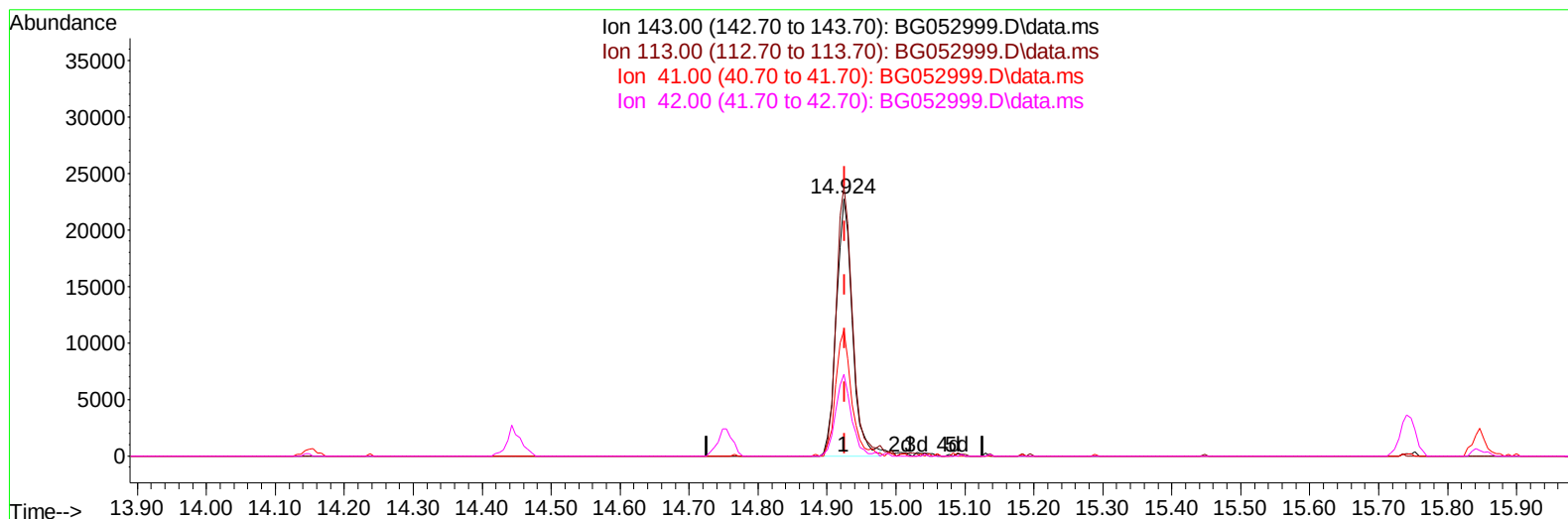
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 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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TIC: BG052999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.924min (-0.001) 28.87 ng/ul m

response 36774

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.82
41.00	47.20	48.82
42.00	29.40	31.75

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	30636	20.000	ng/ul	0.00
20) Naphthalene-d8	10.942	136	127080	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.754	164	98623	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.497	188	234514	20.000	ng/ul	0.00
79) Chrysene-d12	21.780	240	244303	20.000	ng/ul	# 0.00
88) Perylene-d12	25.087	264	238100	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.552	96	4875m	5.621	ng/uL	0.00
4) Pyridine-d5	3.963	84	59161	27.323	ng/ul	0.00
7) Phenol-d5	7.276	99	81124	29.565	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	52761	31.483	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	56957	30.570	ng/ul	0.00
15) 4-Methylphenol-d8	8.827	113	60513	30.444	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	30686	33.545	ng/ul	0.00
24) 2-Nitrophenol-d4	10.019	143	33752	34.981	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	62082	29.576	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	85199	30.711	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	243120	32.088	ng/ul	0.00
49) Acenaphthylene-d8	14.449	160	269059	29.224	ng/ul	0.00
54) 4-Nitrophenol-d4	14.924	143	36774m	28.868	ng/ul	0.00
60) Fluorene-d10	15.741	176	205434	30.880	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.847	200	37749	31.630	ng/ul	0.00
73) Anthracene-d10	17.591	188	346225	31.704	ng/ul	0.00
81) Pyrene-d10	19.877	212	419787	30.716	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.858	264	397843	32.338	ng/ul	0.00

Target Compounds Qvalue

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

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