

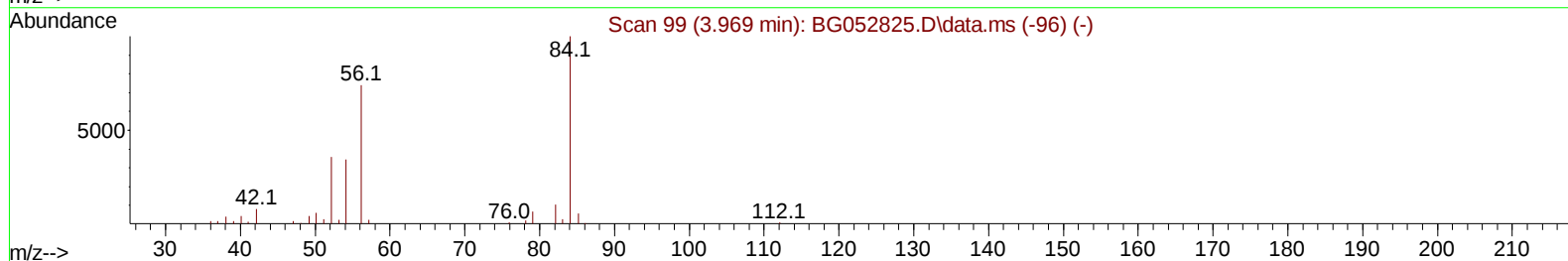
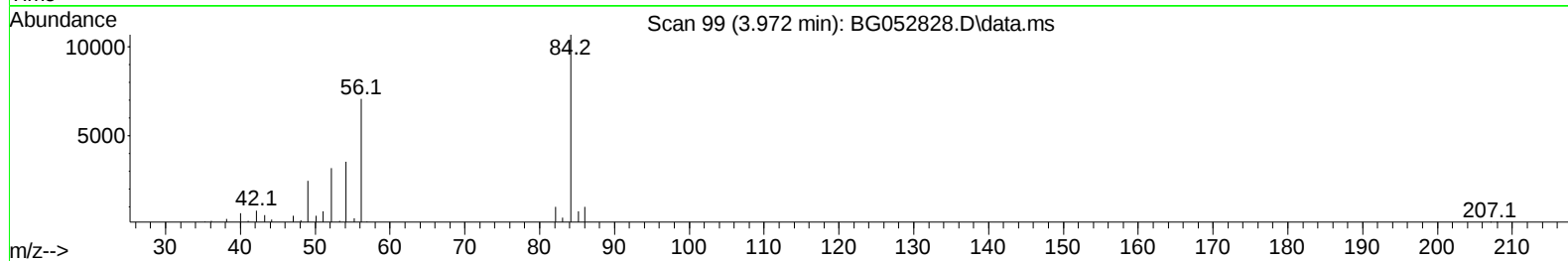
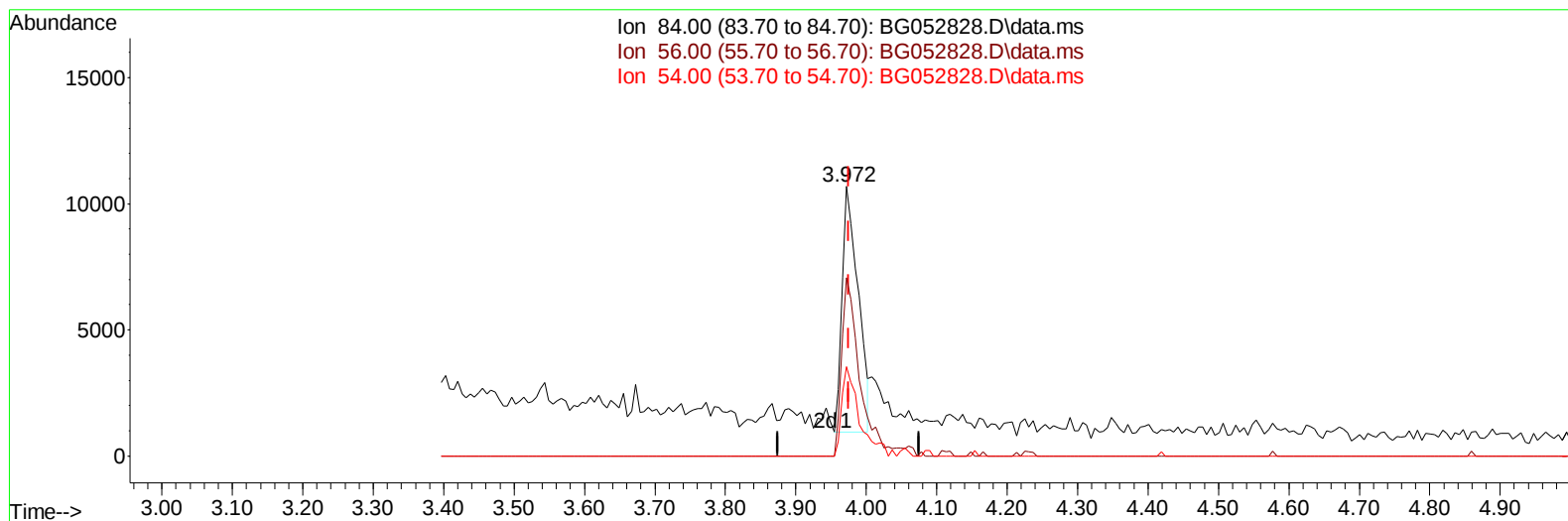
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
 Data File : BG052828.D
 Acq On : 25 Mar 2022 14:07
 Operator : CG/JU
 Sample : N2079-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW618

Manual IntegrationsAPPROVED

Quant Time: Mar 25 15:10:08 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: BG052828.D\data.ms

(4) Pyridine-d5 (S)

3.972min (-0.003) 8.46 ng/u1

response 15301

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	66.24
54.00	28.70	33.20
0.00	0.00	0.00

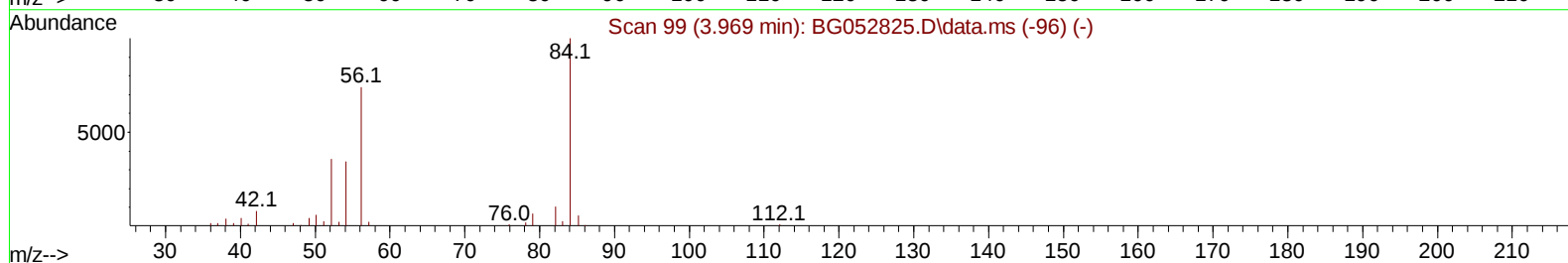
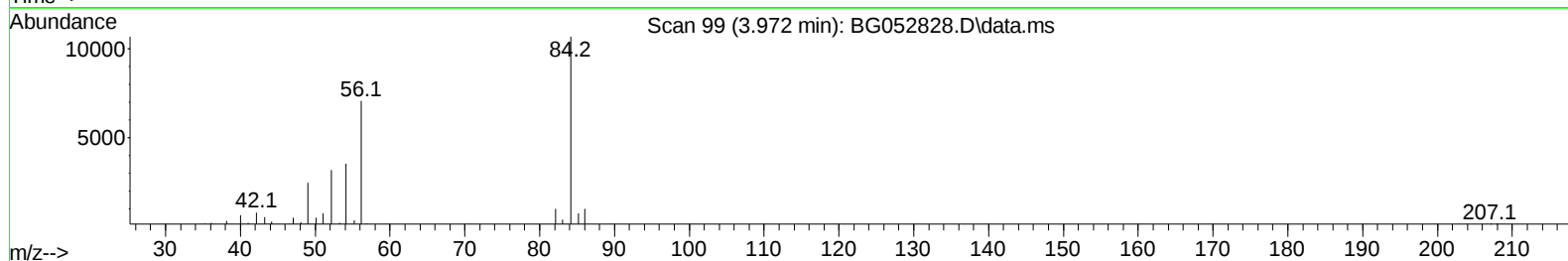
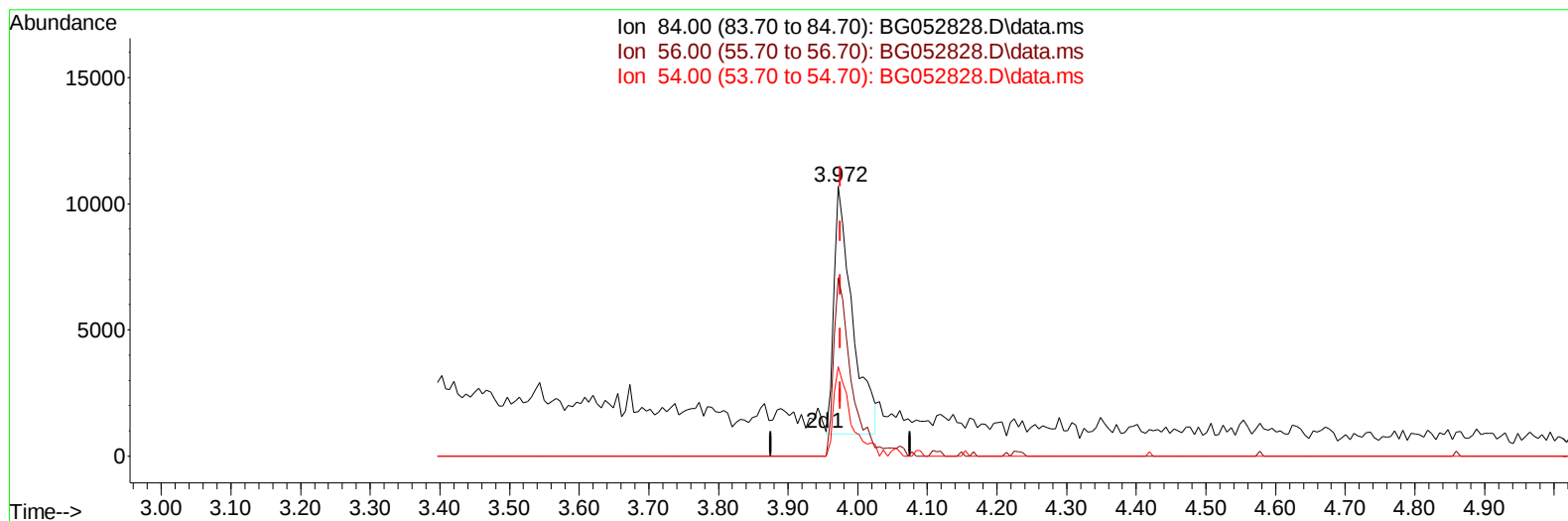
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(4) Pyridine-d5 (S)

3.972min (-0.003) 9.97 ng/ul m

response 18018

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	66.24
54.00	28.70	33.20
0.00	0.00	0.00

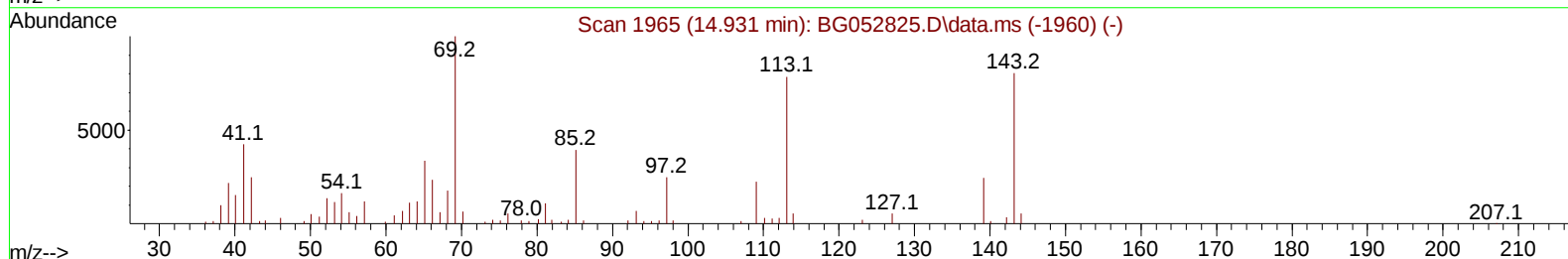
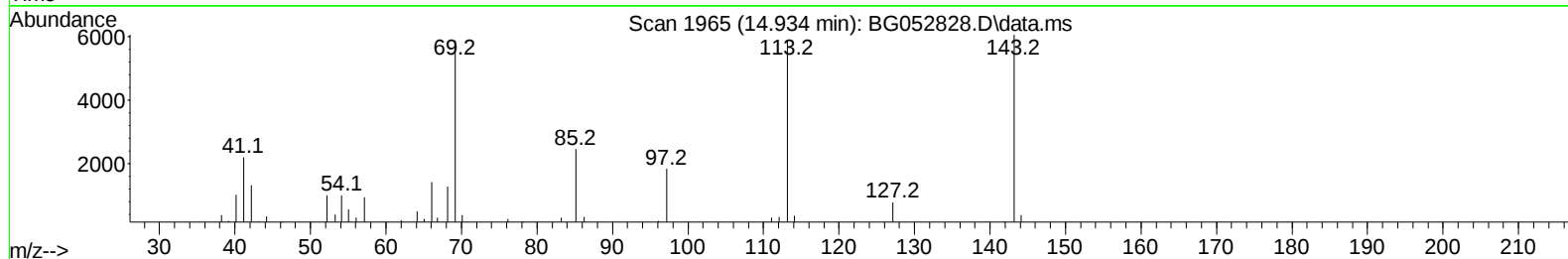
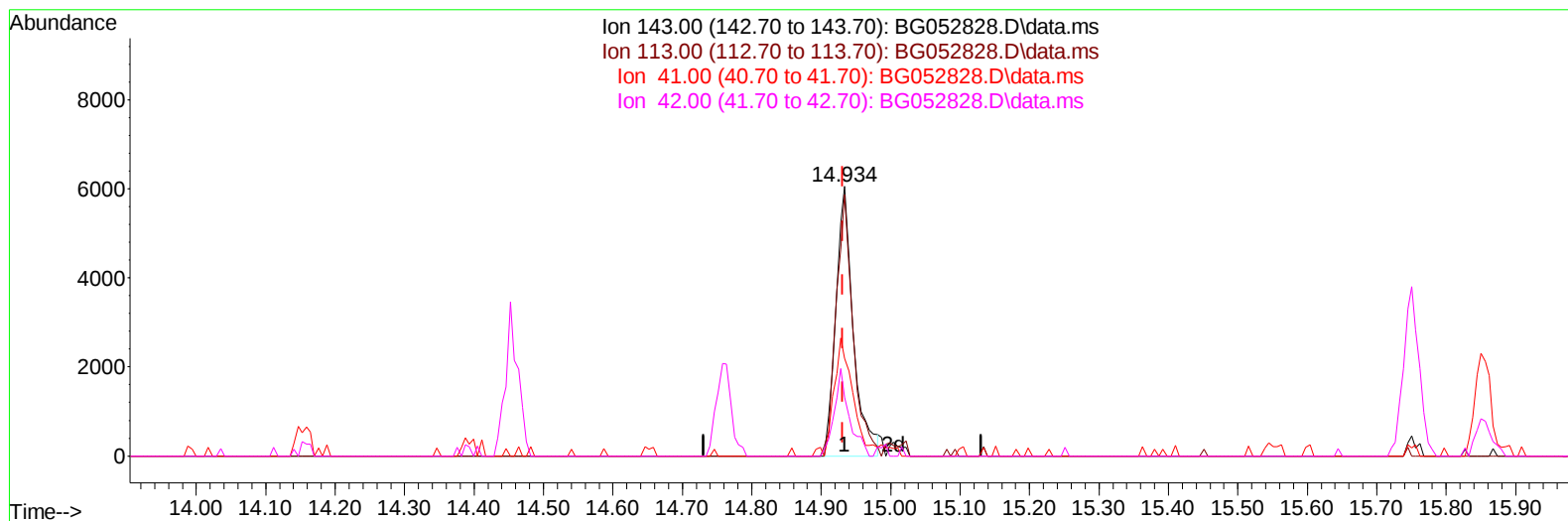
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TIC: BG052828.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.003) 8.68 ng/u1

response 10600

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.09
41.00	47.20	36.26#
42.00	29.40	21.91#

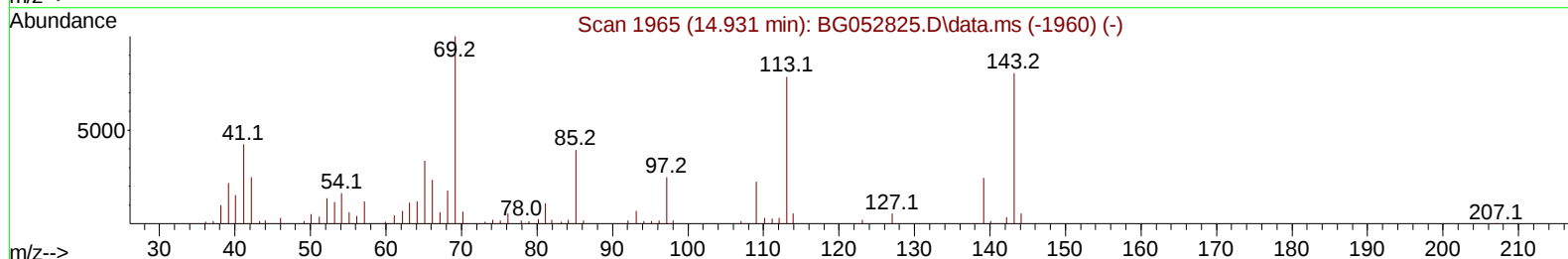
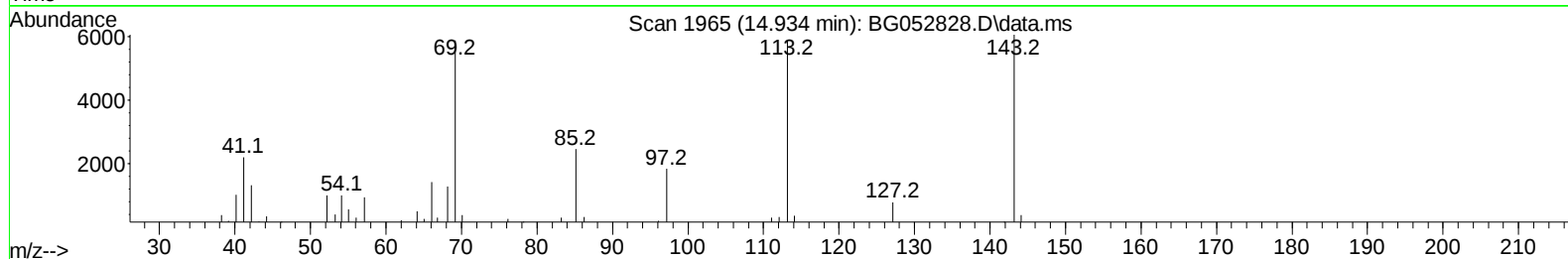
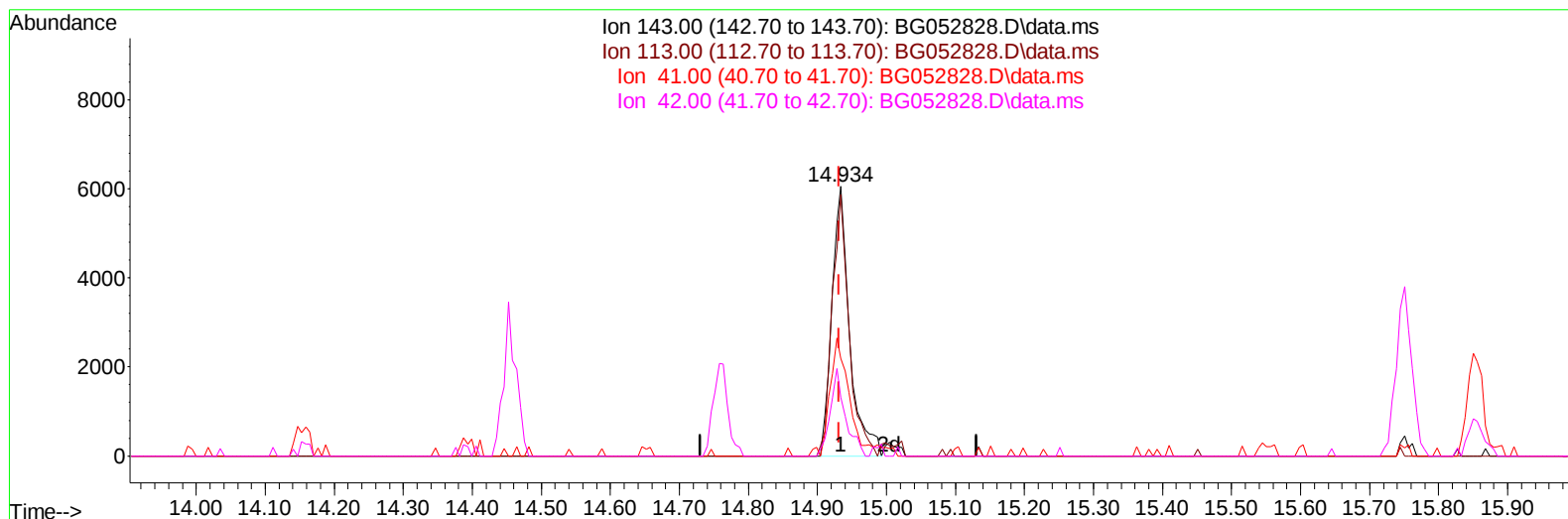
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Operator : CG/JU
Sample : N2079-10
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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TIC: BG052828.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.003) 8.80 ng/ul m

response 10749

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.09
41.00	47.20	36.26#
42.00	29.40	21.91#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.143	152	25576	20.000	ng/ul	0.00
20) Naphthalene-d8	10.951	136	113673	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.764	164	94570	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.501	188	240351	20.000	ng/ul	0.00
79) Chrysene-d12	21.789	240	252924	20.000	ng/ul	0.00
88) Perylene-d12	25.108	264	251832	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.555	96	2815	3.888	ng/uL	0.00
4) Pyridine-d5	3.972	84	18018m	9.968	ng/ul	0.00
7) Phenol-d5	7.285	99	19425	8.480	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.456	67	51313	36.677	ng/ul	0.00
11) 2-Chlorophenol-d4	7.673	132	42579	27.375	ng/ul	0.00
15) 4-Methylphenol-d8	8.836	113	33305	20.071	ng/ul	0.00
21) Nitrobenzene-d5	9.300	128	30744	37.572	ng/ul	0.00
24) 2-Nitrophenol-d4	10.029	143	33889	39.266	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	60128	32.024	ng/ul	0.00
31) 4-Chloroaniline-d4	11.080	131	71287	28.727	ng/ul	0.00
46) Dimethylphthalate-d6	14.158	166	266570	36.691	ng/ul	0.00
49) Acenaphthylene-d8	14.458	160	304994	34.547	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	10749m	8.800	ng/ul	0.00
60) Fluorene-d10	15.750	176	238005	37.309	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.856	200	47139	38.539	ng/ul	0.00
73) Anthracene-d10	17.607	188	425651	38.031	ng/ul	0.00
81) Pyrene-d10	19.886	212	511237	36.132	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.885	264	491765	37.793	ng/ul	0.00

Target Compounds Qvalue

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

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