Quantitation Report (Qedit)

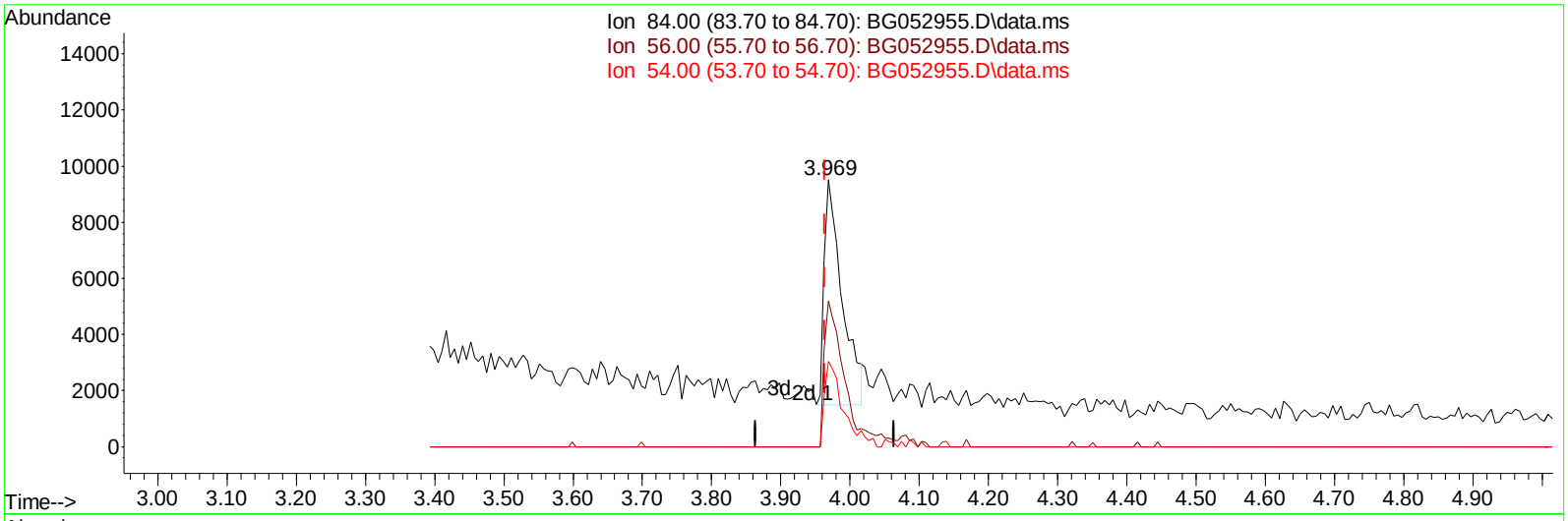
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052955.D
Acq On : 30 Mar 2022 18:43
Operator : CG/JU
Sample : N2068-14
Misc :
ALS Vial : 12 Sample Multiplier: 1

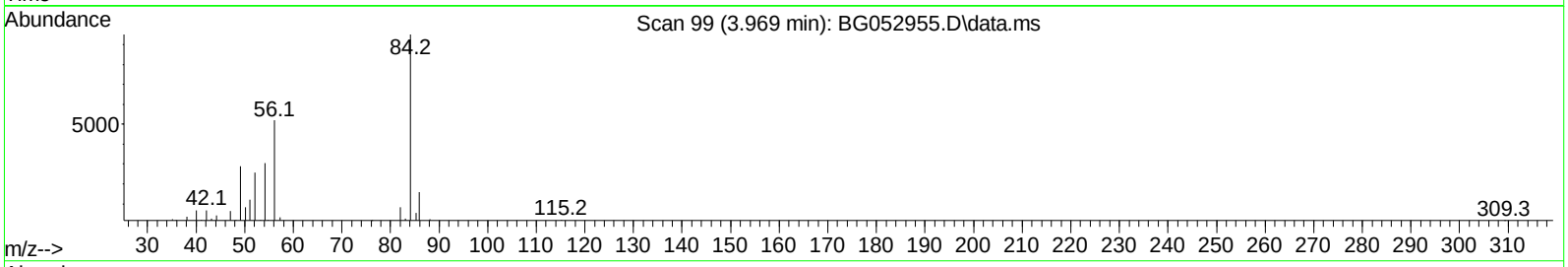
Quant Time: Mar 31 00:53:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

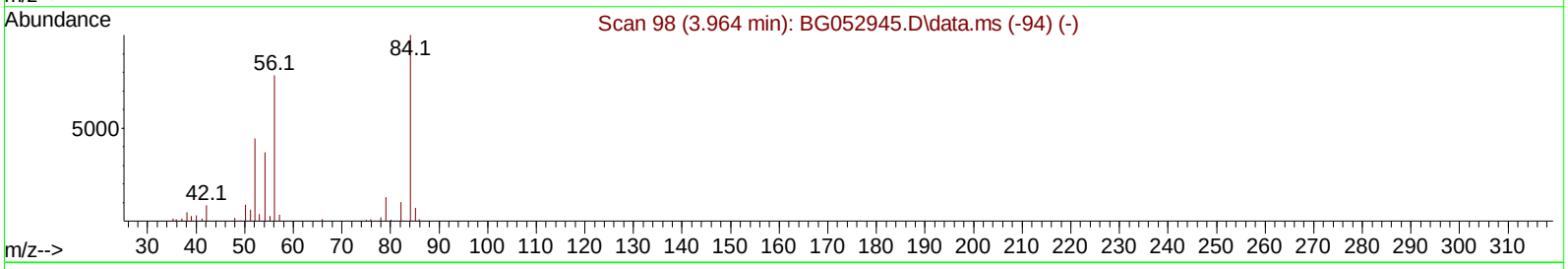
Instrument :
BNA_G
ClientSampleId :
YAZ61

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/06/2022







TIC: BG052955.D\data.ms

(4) Pyridine-d5 (S)

3.969min (+ 0.005) 8.10 ng/ul

response 14342

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.61
54.00	28.70	31.95
0.00	0.00	0.00

Quantitation Report (Qedit)

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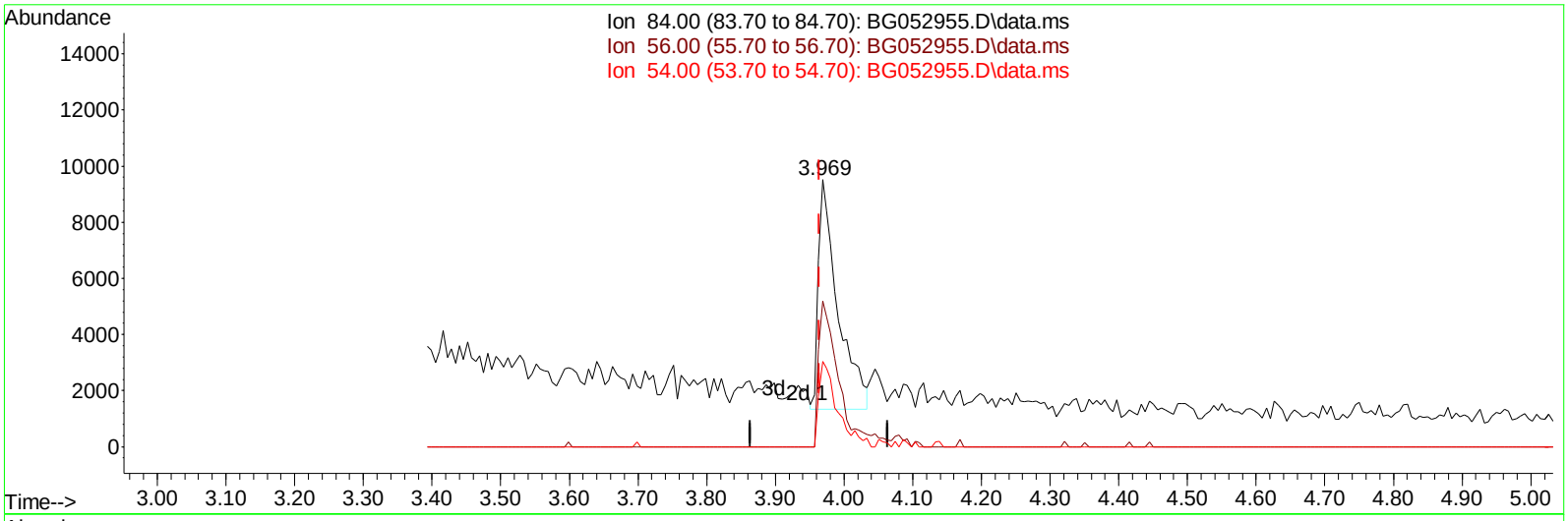

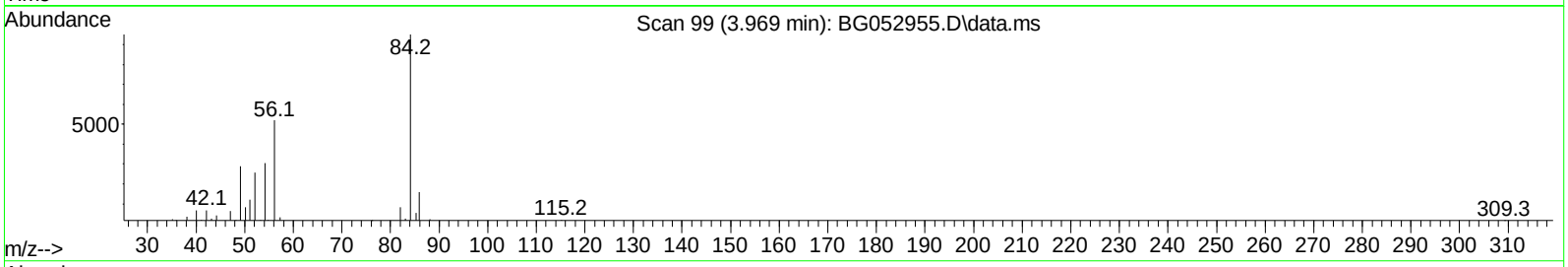

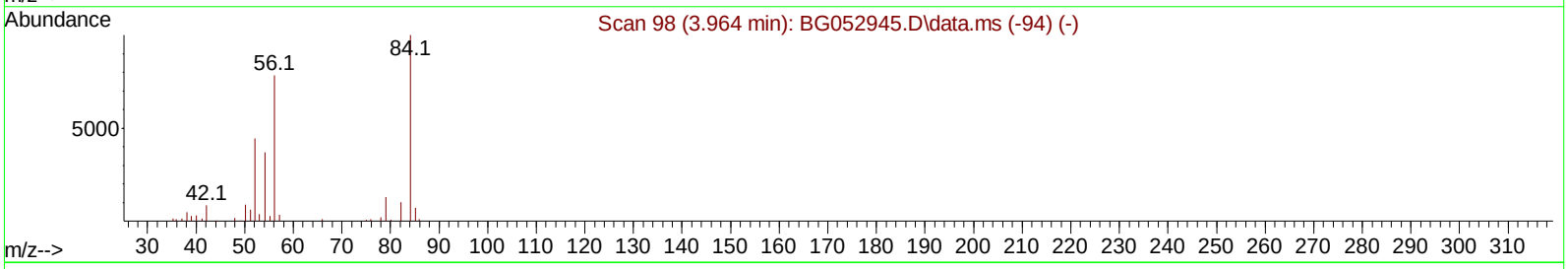

TIC: BG052955.D\data.ms

(4) Pyridine-d5 (S)

3.969min (+ 0.005) 9.10 ng/ul m

response 16103

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.61
54.00	28.70	31.95
0.00	0.00	0.00

Quantitation Report (Qedit)

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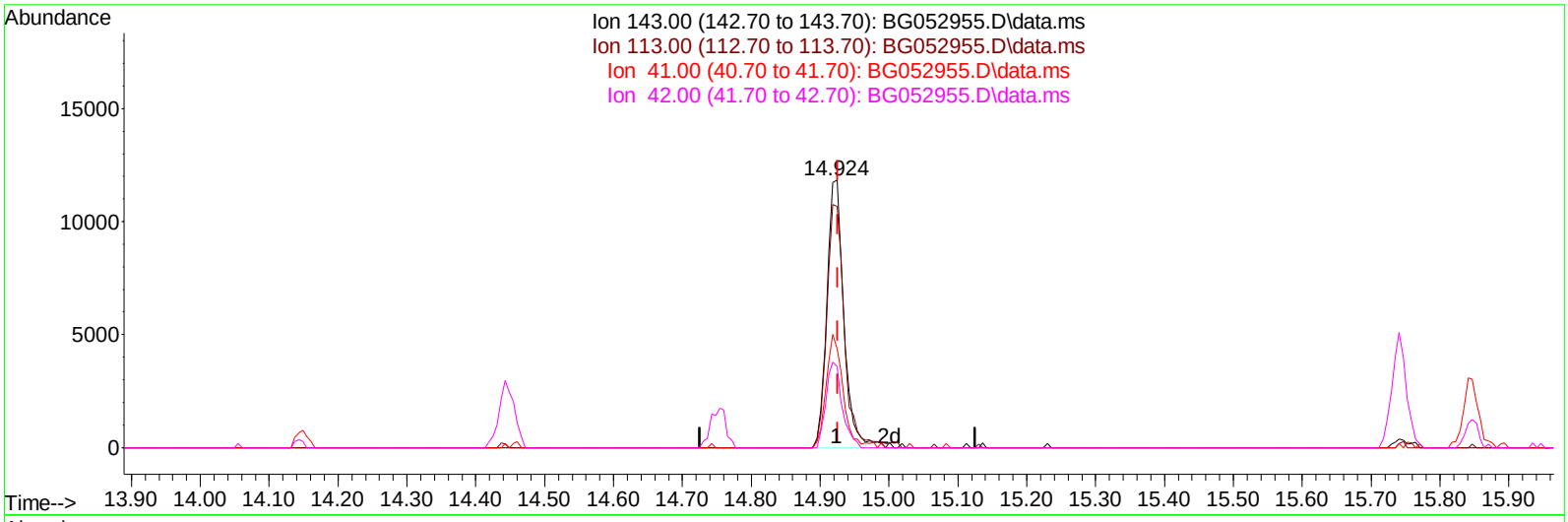

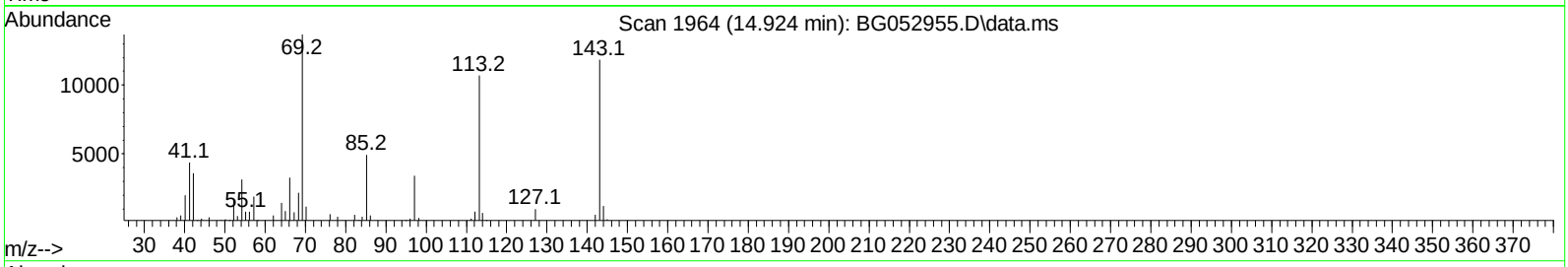

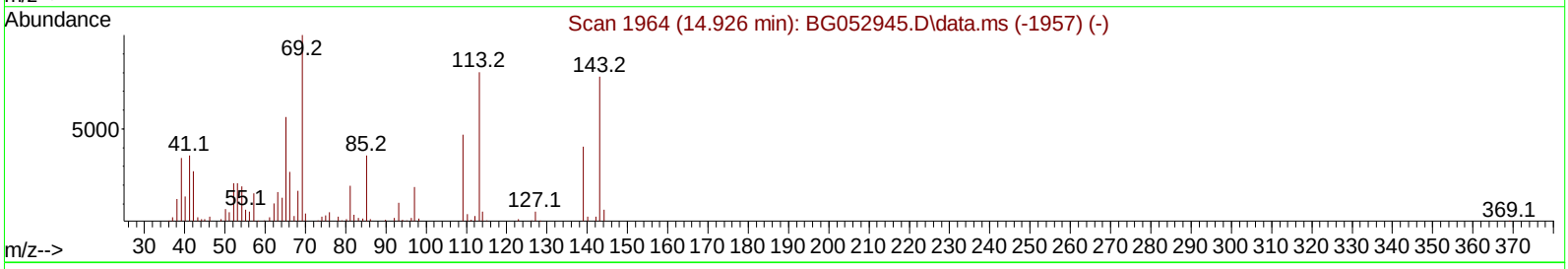

TIC: BG052955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.924min (-0.001) 18.21 ng/ul

response 19945

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	90.30
41.00	47.20	36.88#
42.00	29.40	30.20

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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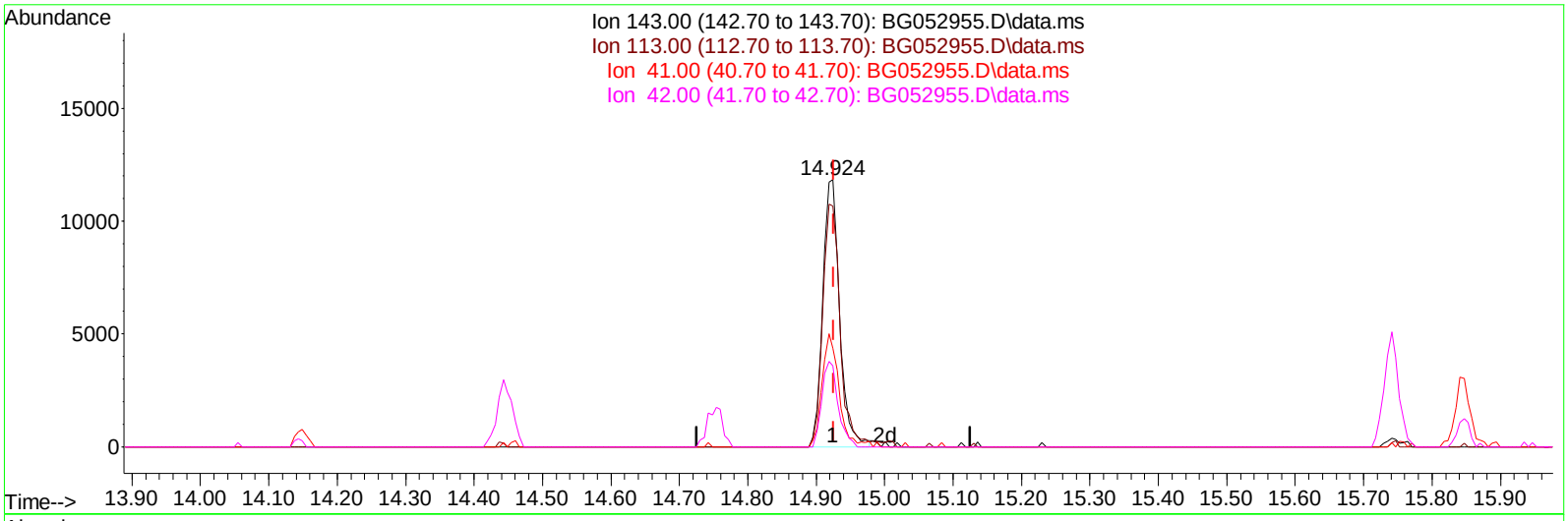

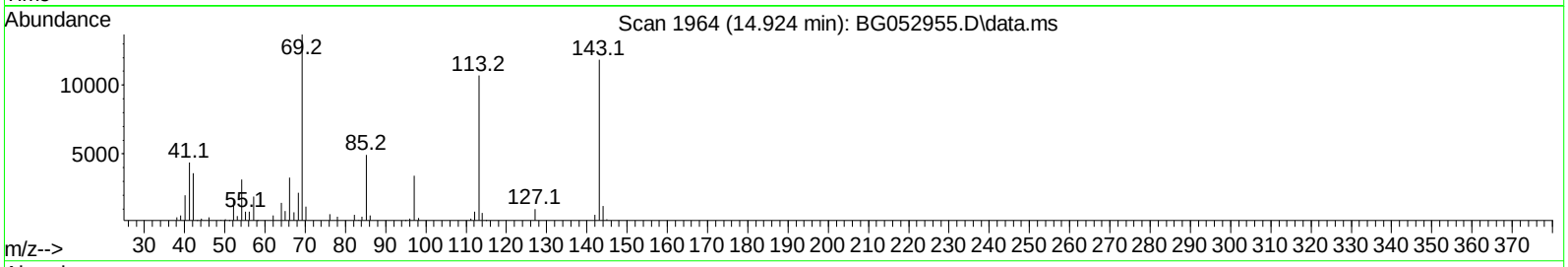

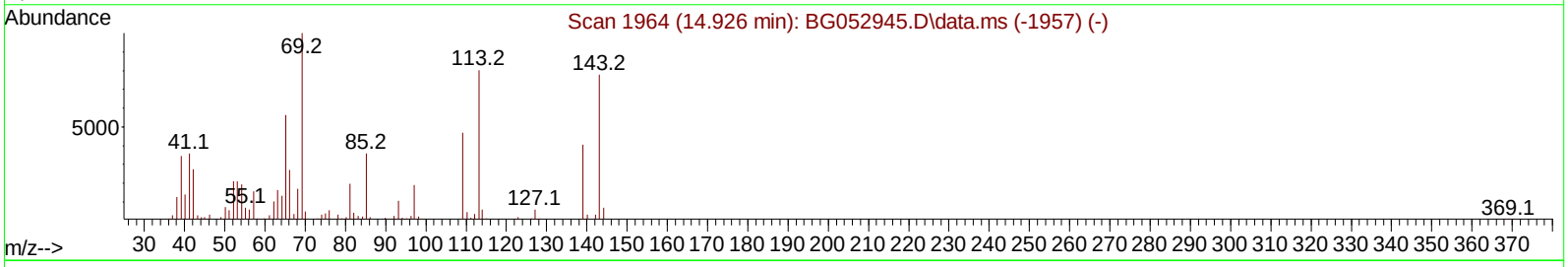

TIC: BG052955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.924min (-0.001) 18.40 ng/ul m

response 20151

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	90.30
41.00	47.20	36.88#
42.00	29.40	30.20

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/06/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	25049	20.000	ng/ul	0.00
20) Naphthalene-d8	10.947	136	109927	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.754	164	84775	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.491	188	209805	20.000	ng/ul	0.00
79) Chrysene-d12	21.780	240	230698	20.000	ng/ul	0.00
88) Perylene-d12	25.087	264	226664	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.546	96	6824	9.623	ng/uL	0.00
4) Pyridine-d5	3.969	84	16103m	9.096	ng/ul	0.00
7) Phenol-d5	7.282	99	35713	15.918	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	62162	45.366	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	53861	35.356	ng/ul	0.00
15) 4-Methylphenol-d8	8.827	113	44007	27.078	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	34809	43.990	ng/ul	0.00
24) 2-Nitrophenol-d4	10.019	143	40402	48.408	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	68256	37.592	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	58563	24.404	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	306911	47.124	ng/ul	0.00
49) Acenaphthylene-d8	14.449	160	323272	40.848	ng/ul	0.00
54) 4-Nitrophenol-d4	14.924	143	20151m	18.403	ng/ul	0.00
60) Fluorene-d10	15.741	176	266168	46.545	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.847	200	56785	53.185	ng/ul	0.00
73) Anthracene-d10	17.591	188	477791	48.905	ng/ul	0.00
81) Pyrene-d10	19.877	212	608453	47.146	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.864	264	610709	52.145	ng/ul	0.00

Target Compounds Qvalue

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

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