

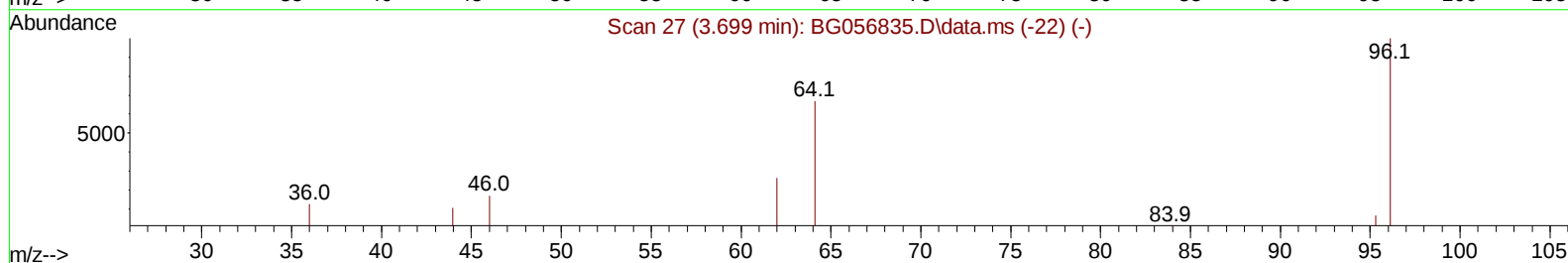
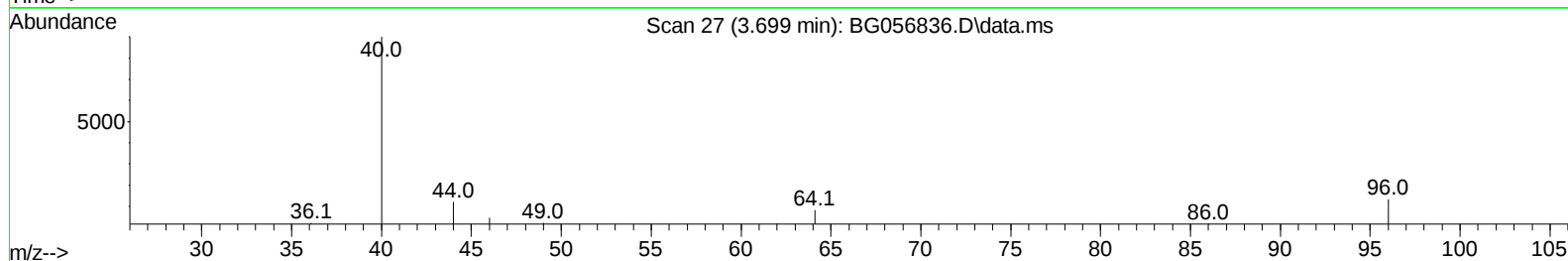
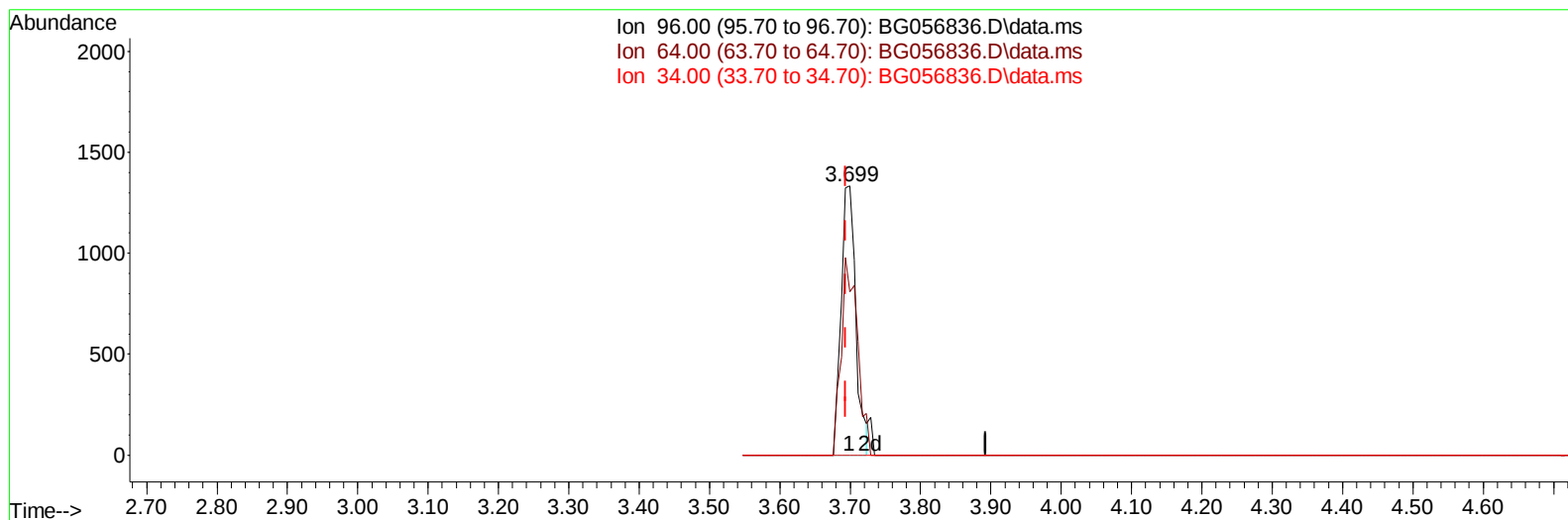
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056836.D
Acq On : 6 Mar 2023 11:07
Operator : CG/JU
Sample : PB151218BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK218

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/07/2023
Supervised By :Jagrut Upadhyay 03/07/2023

Quant Time: Mar 07 01:50:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 02 00:12:28 2023
Response via : Initial Calibration



TIC: BG056836.D\data.ms

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

3.699min (+ 0.006) 7.35 ng/uL

response 1901

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	60.40	60.74
34.00	0.00	0.00
0.00	0.00	0.00

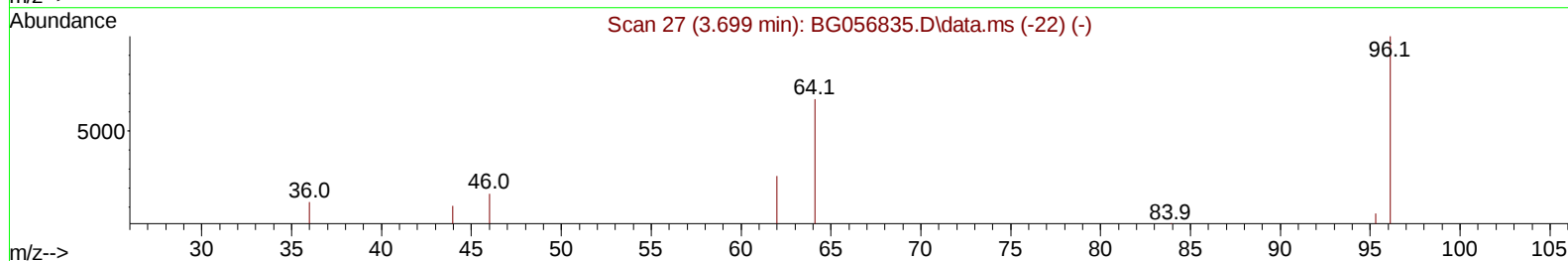
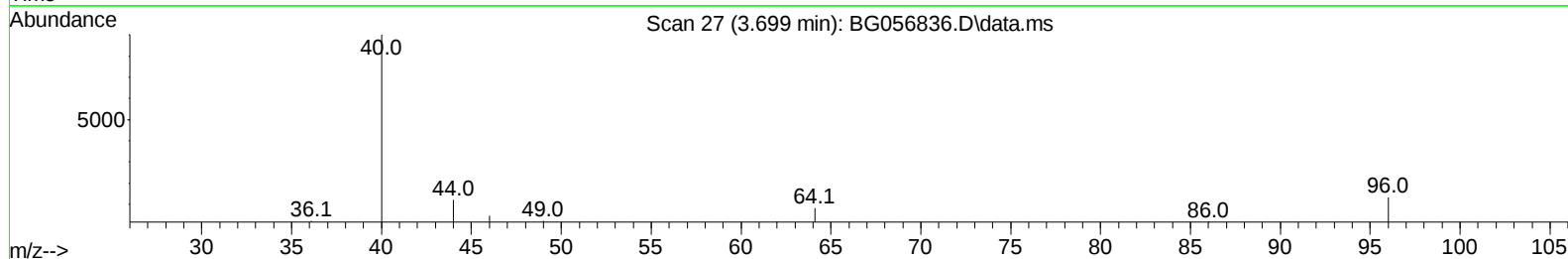
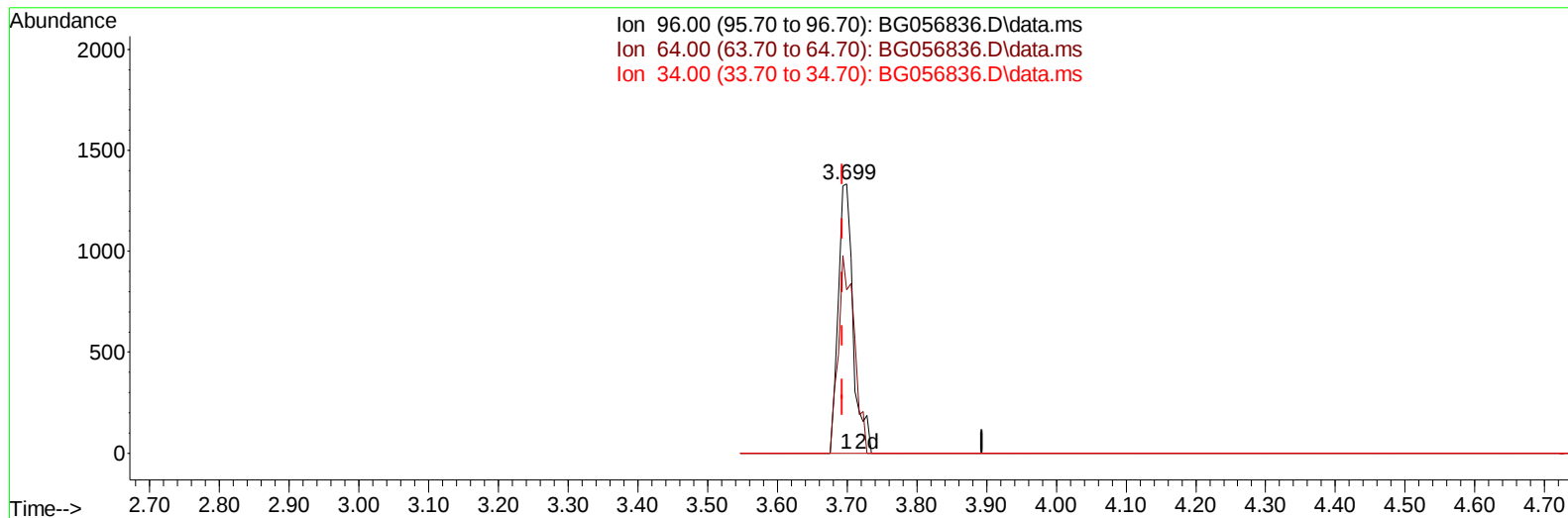
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TIC: BG056836.D\data.ms

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

3.699min (+ 0.006) 7.61 ng/uL m

response 1967

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	60.40	60.74
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :

BNA_G

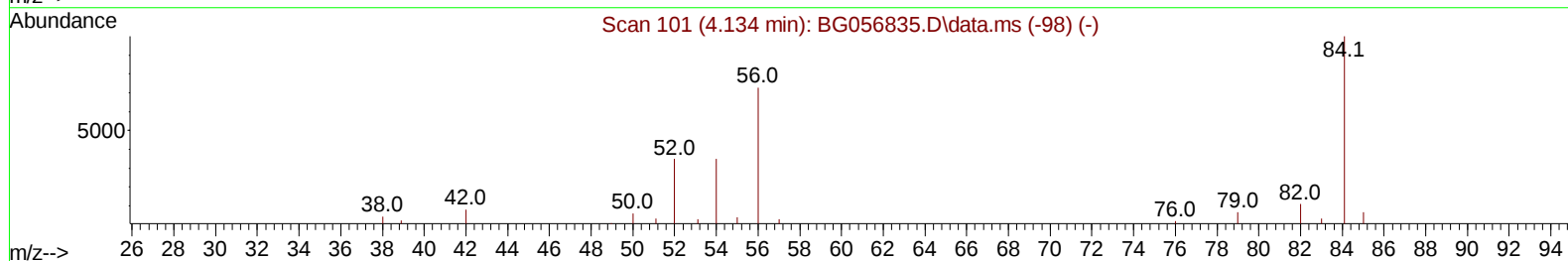
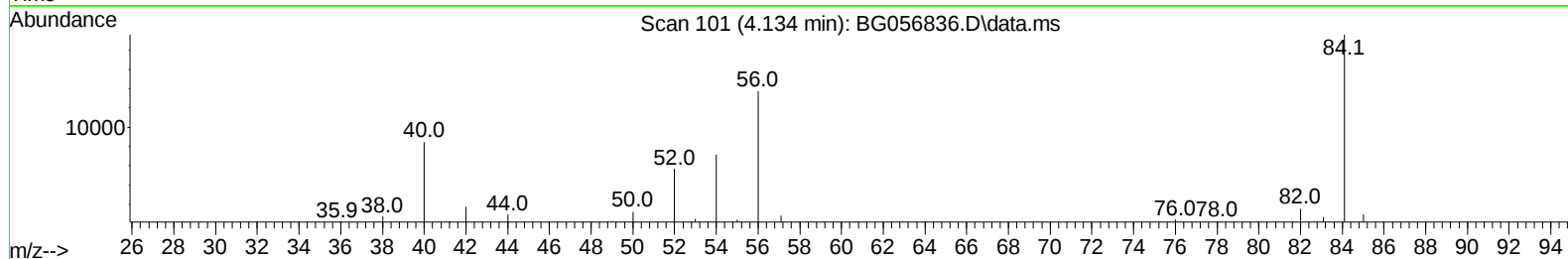
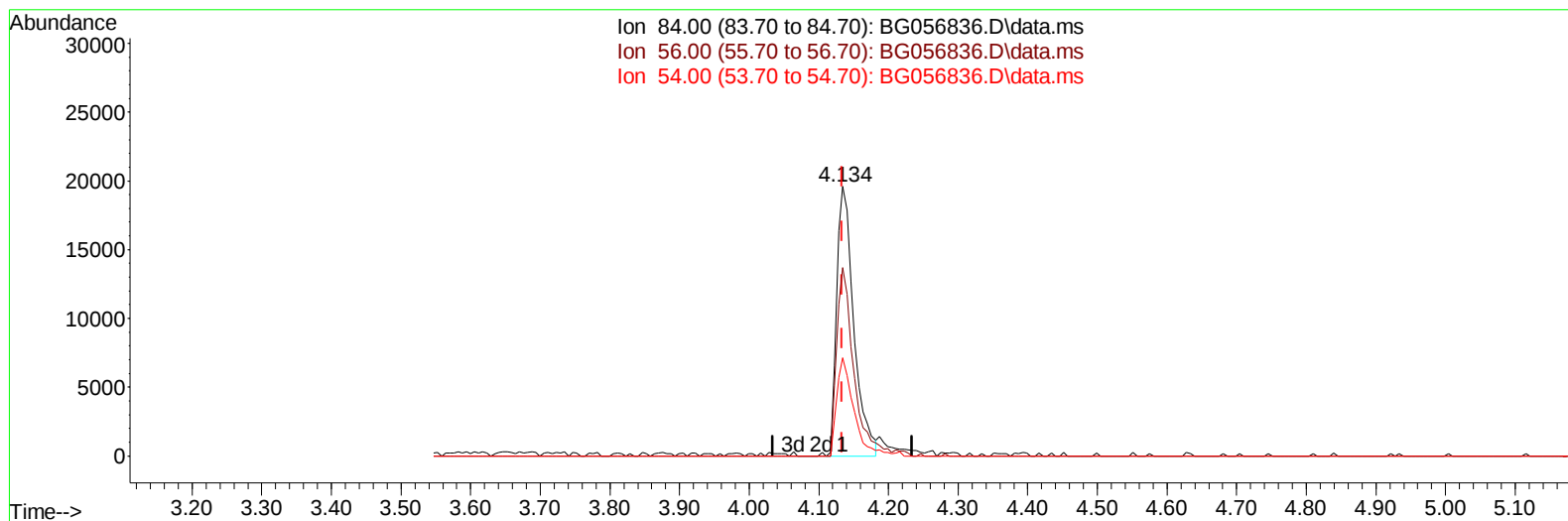
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(4) Pyridine-d5 (S)

4.134min (+ 0.000) 31.99 ng/u1

response 33393

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	67.40	69.91
54.00	38.20	36.54
0.00	0.00	0.00

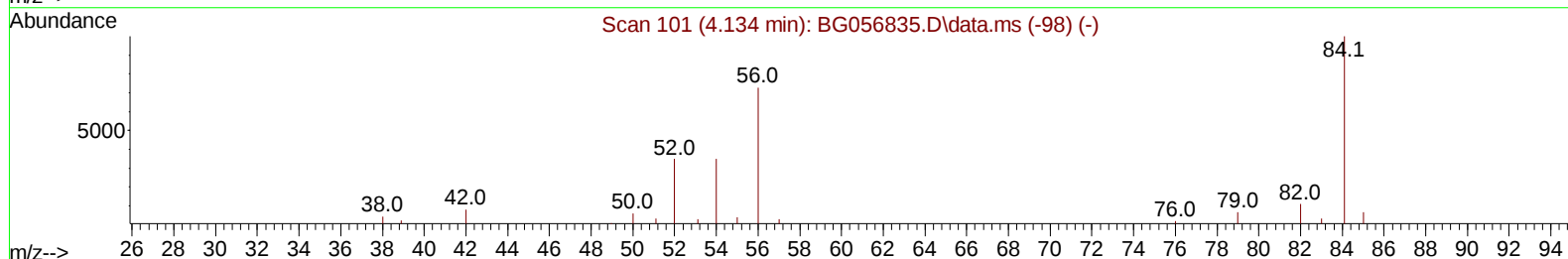
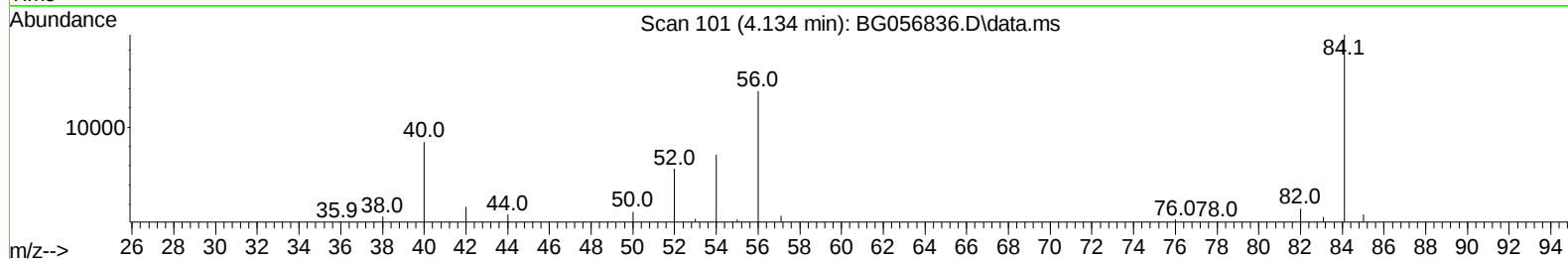
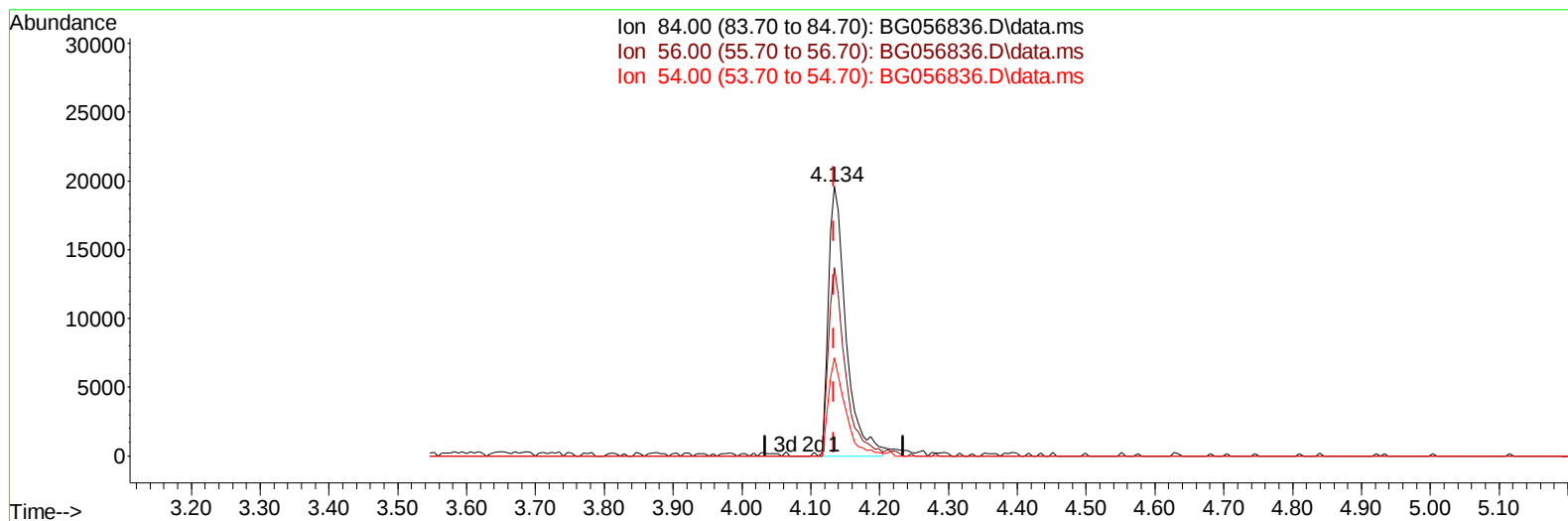
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Instrument :
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TIC: BG056836.D\data.ms

(4) Pyridine-d5 (S)

4.134min (+ 0.000) 33.21 ng/ul m

response 34673

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	67.40	69.91
54.00	38.20	36.54
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.412	152	14347	20.000	ng/ul	0.00
20) Naphthalene-d8	11.249	136	65405	20.000	ng/ul	-0.01
38) Acenaphthene-d10	15.016	164	54177	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.748	188	135986	20.000	ng/ul	0.00
79) Chrysene-d12	22.054	240	133747	20.000	ng/ul	-0.01
88) Perylene-d12	25.574	264	141486	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.699	96	1967m	7.605	ng/uL	0.00
4) Pyridine-d5	4.134	84	34673m	33.213	ng/ul	0.00
7) Phenol-d5	7.542	99	36880	25.688	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.718	67	24169	27.598	ng/ul	0.00
11) 2-Chlorophenol-d4	7.936	132	24033	25.207	ng/ul	0.00
15) 4-Methylphenol-d8	9.105	113	27235	24.827	ng/ul	0.00
21) Nitrobenzene-d5	9.587	128	13776	25.133	ng/ul	0.00
24) 2-Nitrophenol-d4	10.315	143	15026	23.030	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.856	165	30439	24.764	ng/ul	0.00
31) 4-Chloroaniline-d4	11.373	131	40942	25.538	ng/ul	0.00
46) Dimethylphthalate-d6	14.405	166	108531	24.829	ng/ul	0.00
49) Acenaphthylene-d8	14.716	160	128978	25.962	ng/ul	0.00
54) 4-Nitrophenol-d4	15.174	143	15302	20.304	ng/ul	0.00
60) Fluorene-d10	15.997	176	102031	25.234	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.097	200	17055	18.746	ng/ul	0.00
73) Anthracene-d10	17.848	188	173809	26.346	ng/ul	0.00
81) Pyrene-d10	20.110	212	208464	23.034	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.333	264	201604	26.399	ng/ul	-0.02

Target Compounds Qvalue

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

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