

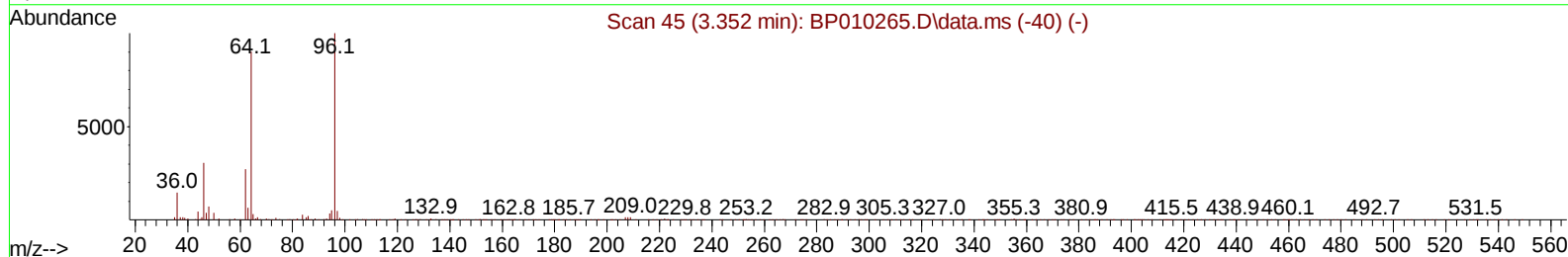
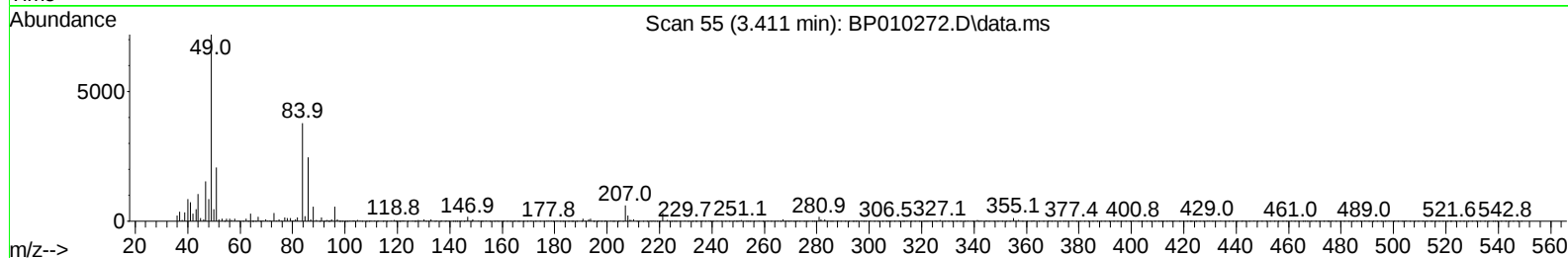
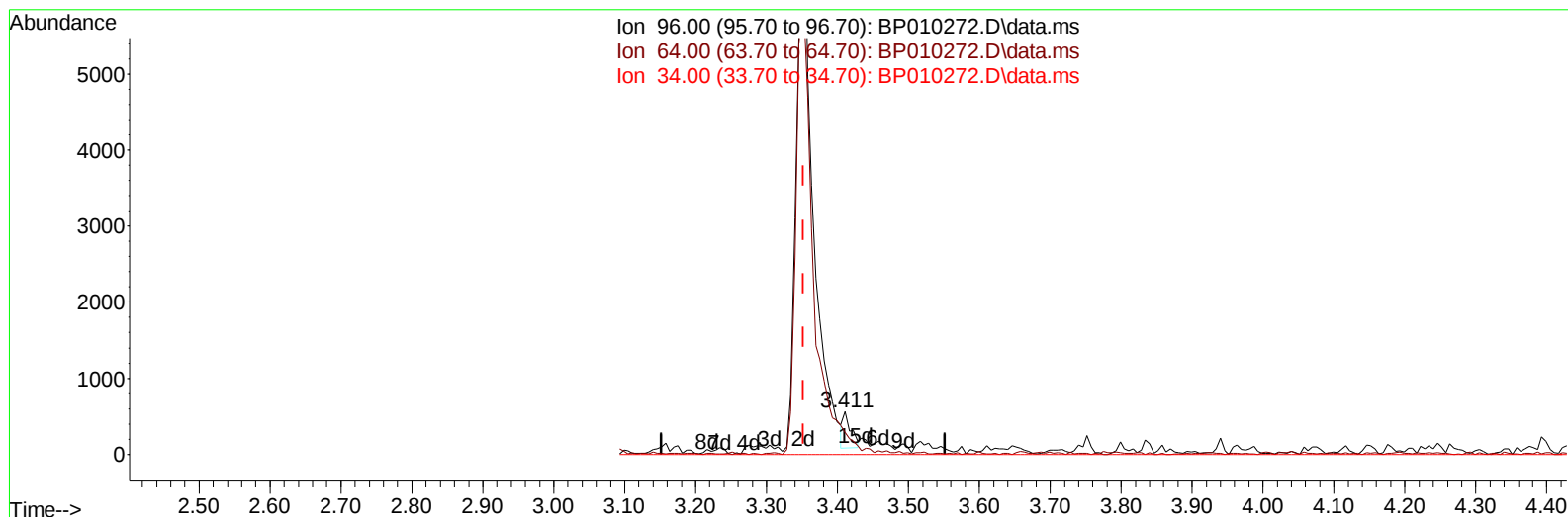
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
 Data File : BP010272.D
 Acq On : 10 May 2022 19:06
 Operator : CG/JU
 Sample : N2678-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0C2

Manual IntegrationsAPPROVED

Quant Time: May 11 01:18:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 10 15:59:46 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/11/2022
 Supervised By :mohammad ahmed 05/12/2022



TIC: BP010272.D\data.ms

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

3.411min (+ 0.059) 0.13 ng/uL

response 353

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	53.04
34.00	0.00	0.00
0.00	0.00	0.00

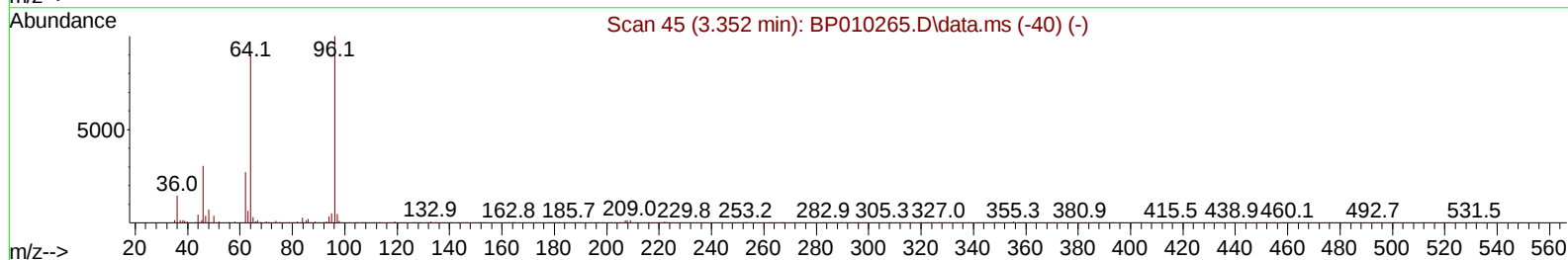
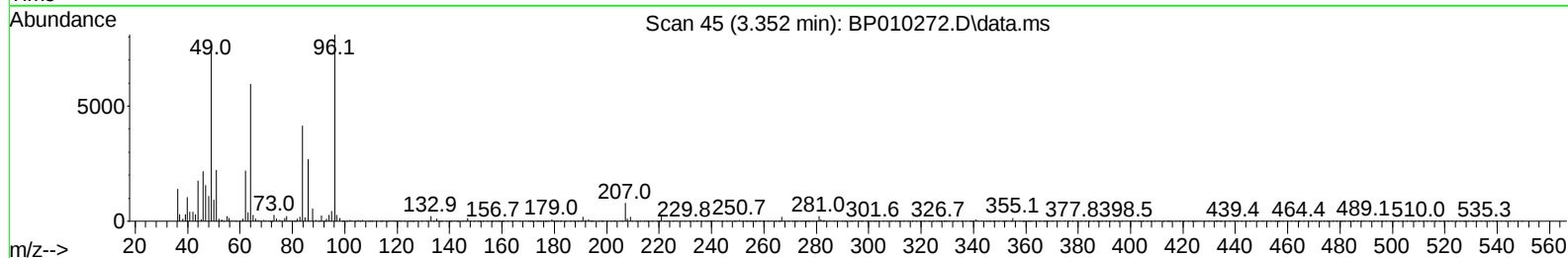
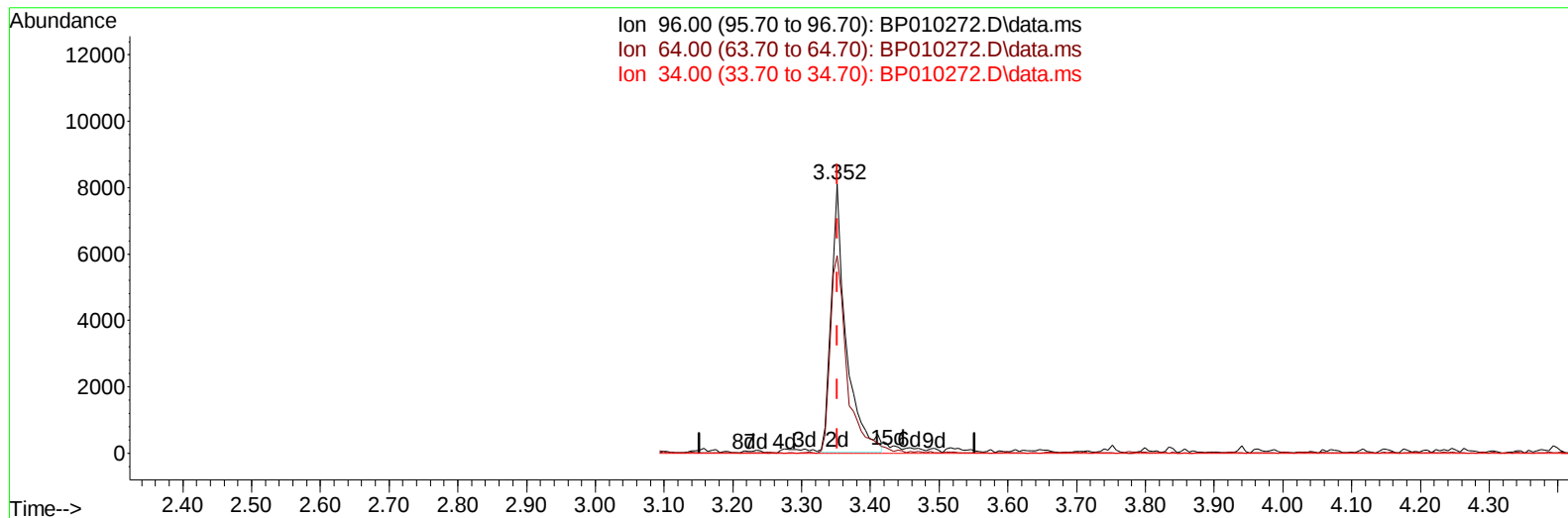
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TIC: BP010272.D\data.ms

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

3.352min (-0.000) 4.61 ng/uL m

response 12165

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.41#
34.00	0.00	0.00
0.00	0.00	0.00

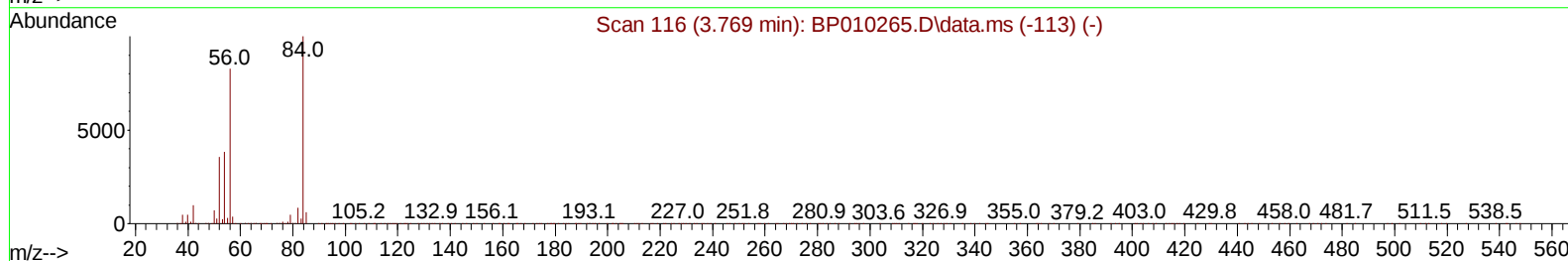
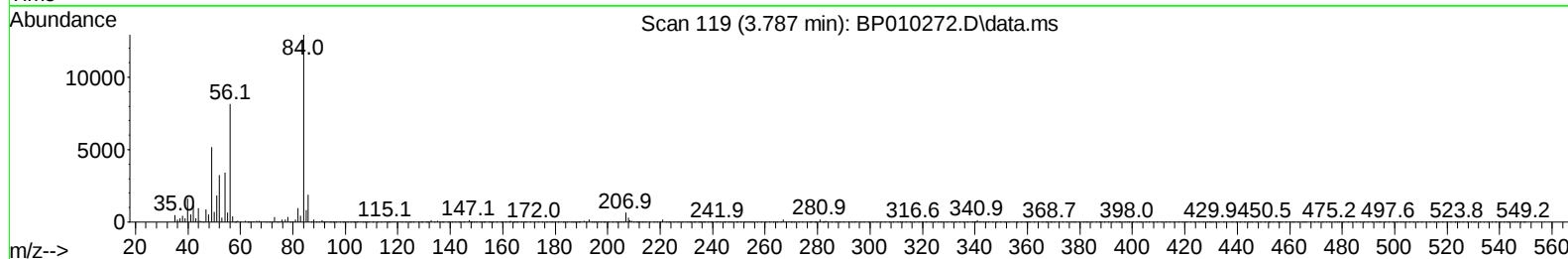
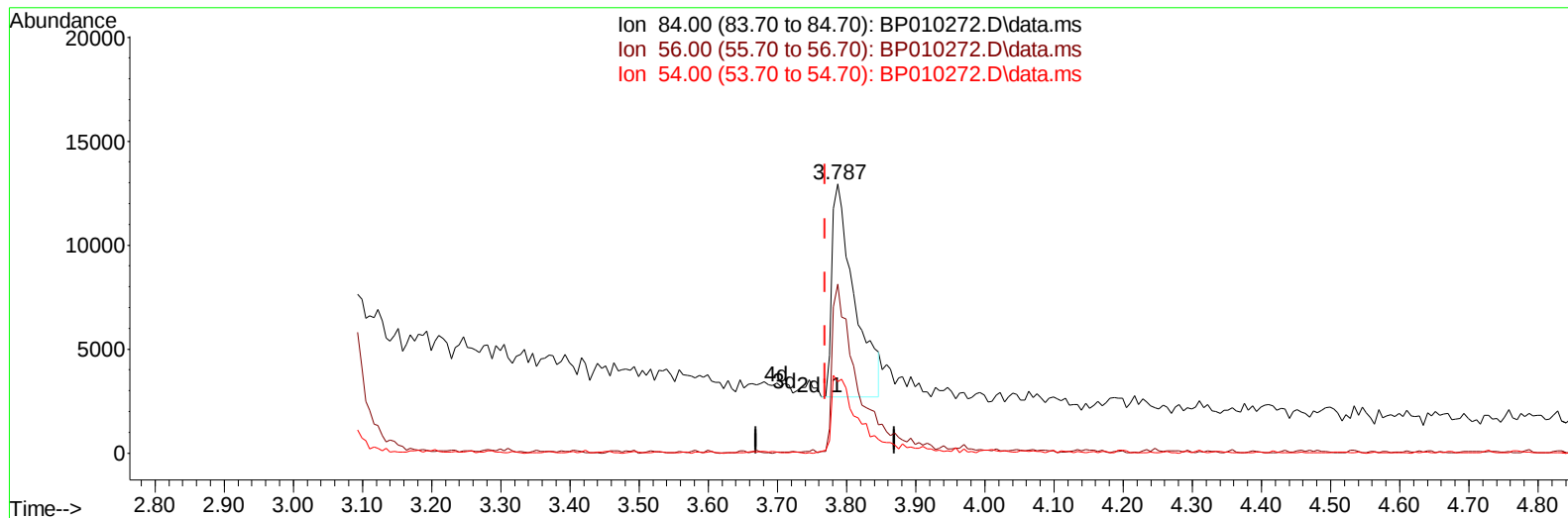
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP010272.D\data.ms

(4) Pyridine-d5 (S)

3.787min (+ 0.018) 3.12 ng/u1

response 22899

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	62.93#
54.00	24.80	26.44
0.00	0.00	0.00

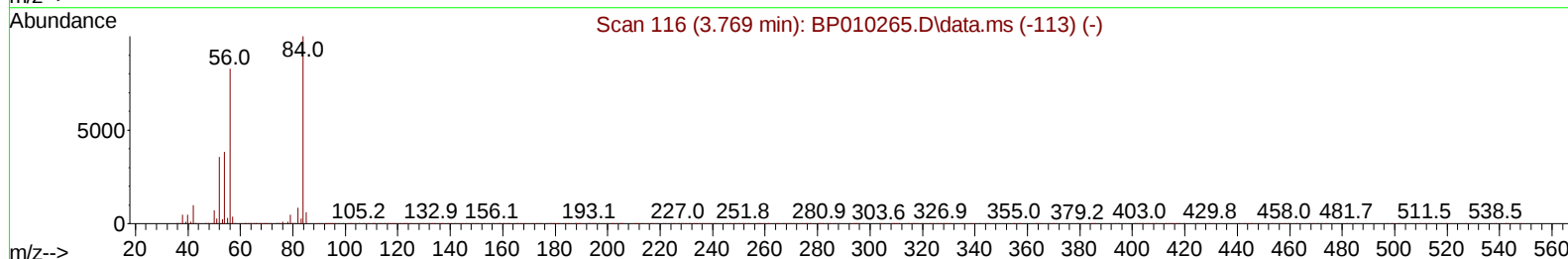
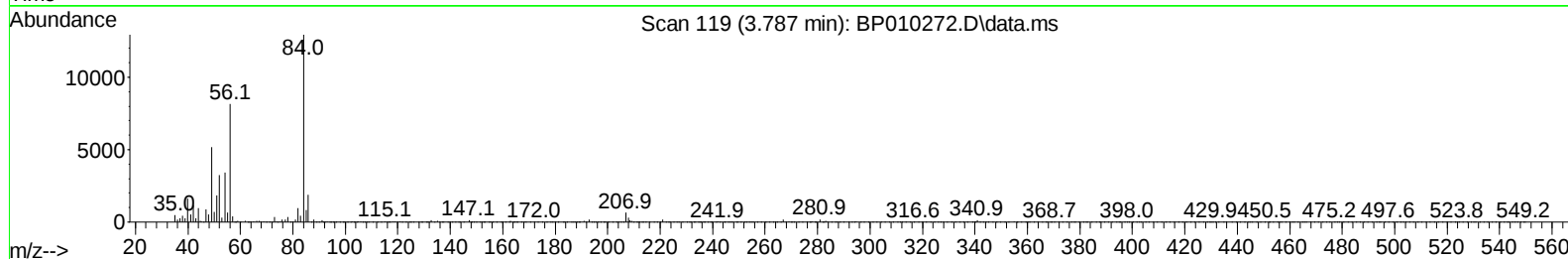
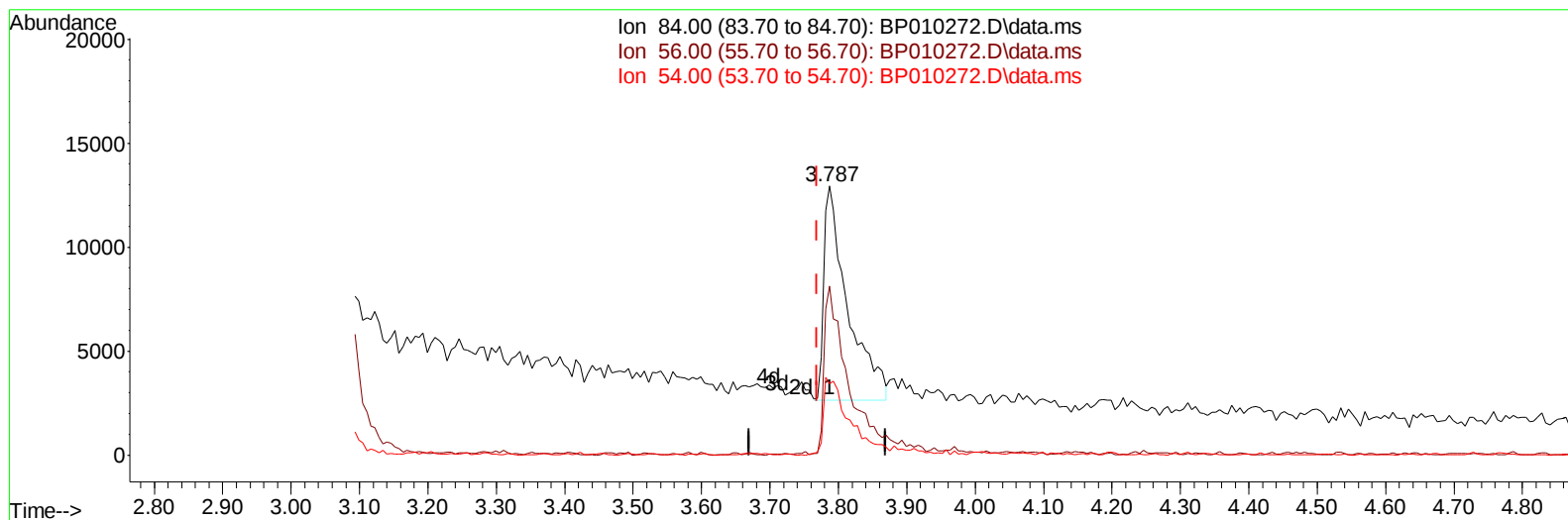
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 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP010272.D\data.ms

(4) Pyridine-d5 (S)

3.787min (+ 0.018) 3.40 ng/u1 m

response 24969

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	62.93#
54.00	24.80	26.44
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	7.910	152	112169	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	491771	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.539	164	353066	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.292	188	794171	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.374	240	823948	20.000	ng/ul	#-0.01
88) Perylene-d12	23.810	264	865091	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	12165m	4.607	ng/uL	0.00
4) Pyridine-d5	3.787	84	24969m	3.398	ng/ul	0.02
7) Phenol-d5	7.075	99	49935	4.793	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	180754	28.046	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	155919	20.321	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	106873	12.478	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	108375	27.485	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	112448	25.012	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	186810	22.829	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	168762	13.970	ng/ul	-0.01
46) Dimethylphthalate-d6	13.951	166	961823	34.195	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	931822	27.786	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	23937	4.437	ng/ul	0.00
60) Fluorene-d10	15.534	176	804689	33.892	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.651	200	153043	28.141	ng/ul	-0.01
73) Anthracene-d10	17.392	188	1388312	35.698	ng/ul	-0.01
81) Pyrene-d10	19.622	212	1723619	37.855	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.651	264	1720164	37.285	ng/ul	-0.02

Target Compounds Qvalue

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

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