

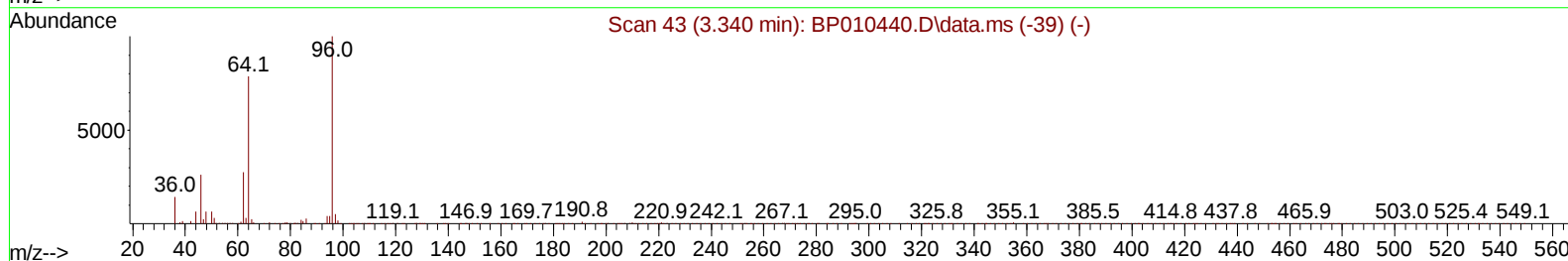
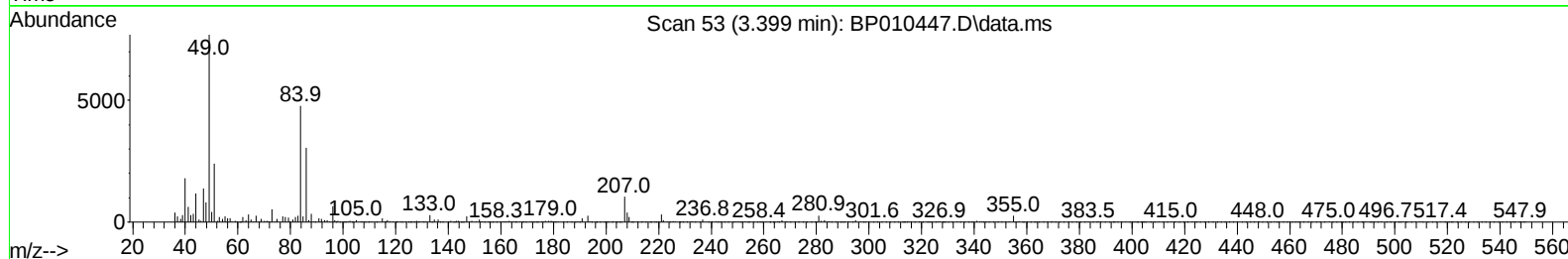
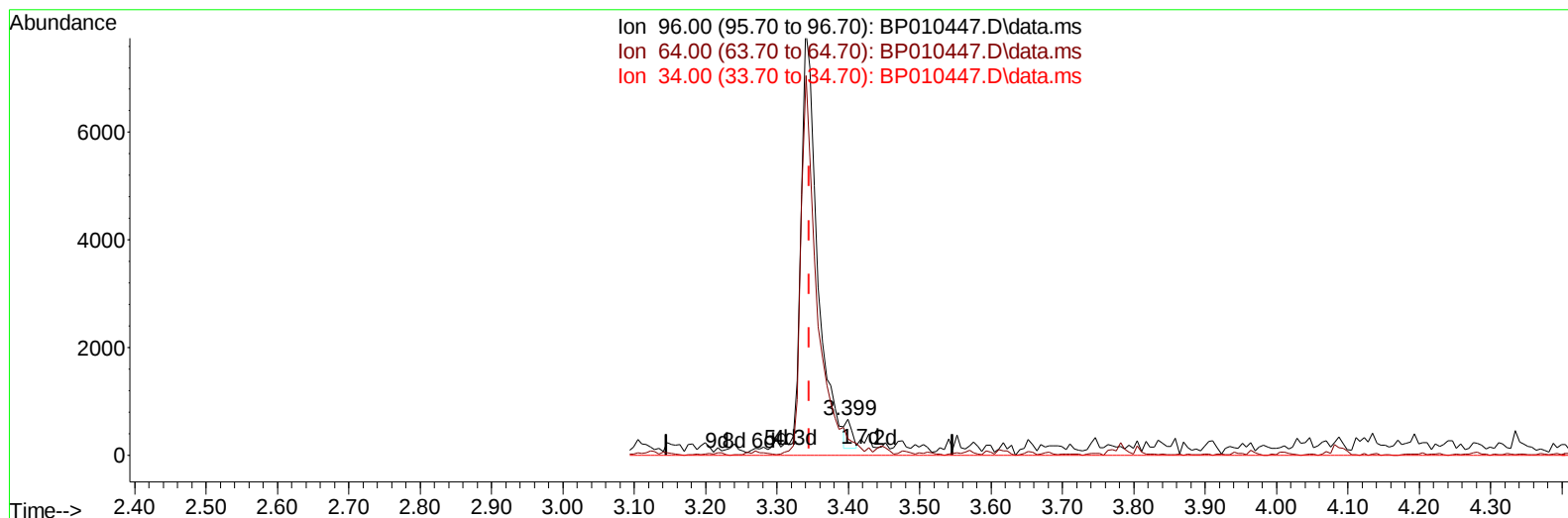
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010447.D
 Acq On : 20 May 2022 21:59
 Operator : CG/JU
 Sample : N2843-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4J1

Manual IntegrationsAPPROVED

Quant Time: May 20 23:04:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/23/2022
 Supervised By :mohammad ahmed 05/25/2022



TIC: BP010447.D\data.ms

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

3.399min (+ 0.053) 0.08 ng/uL

response 322

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.84
34.00	0.00	0.00
0.00	0.00	0.00

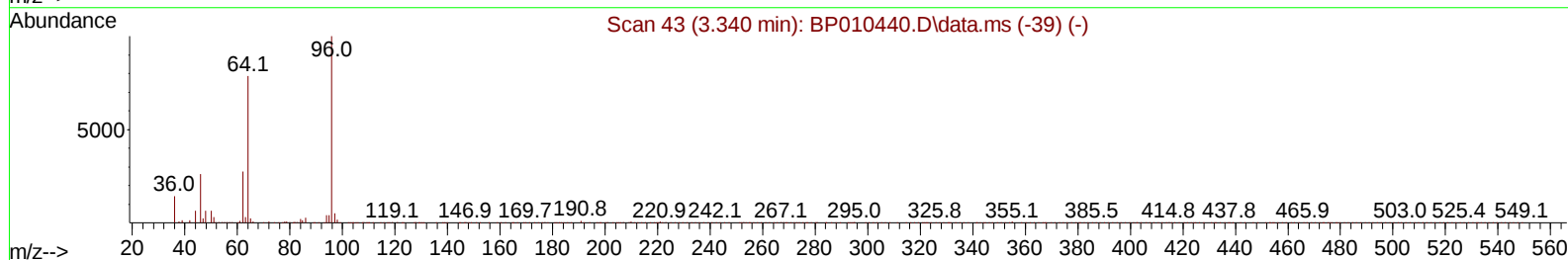
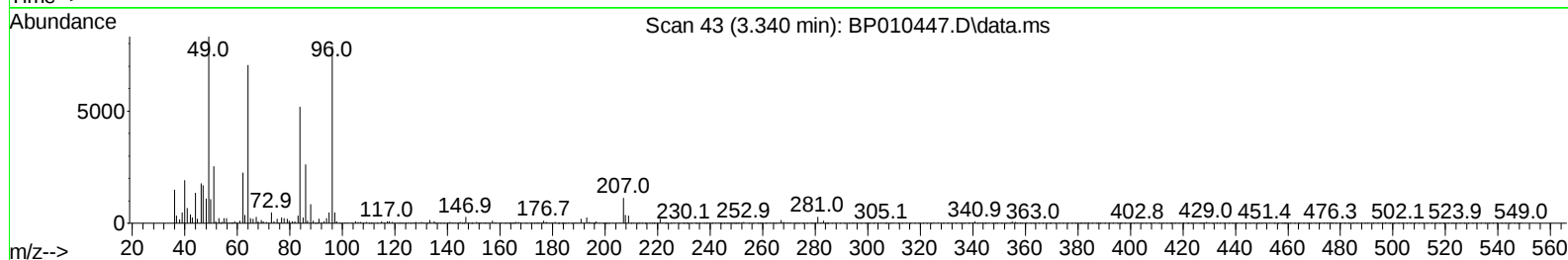
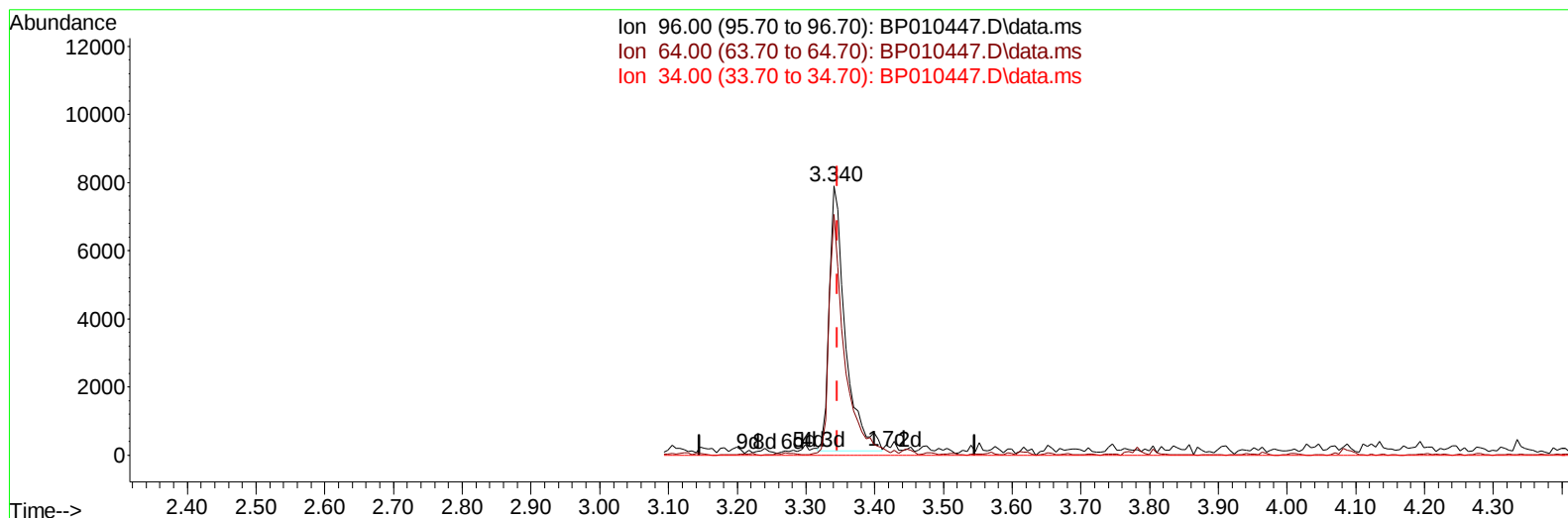
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TIC: BP010447.D\data.ms

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

3.340min (-0.006) 3.19 ng/uL m

response 12597

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.40#
34.00	0.00	0.00
0.00	0.00	0.00

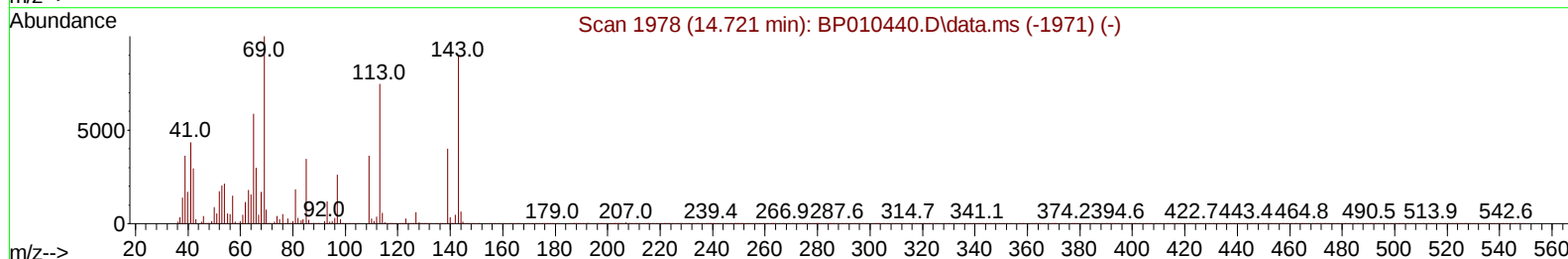
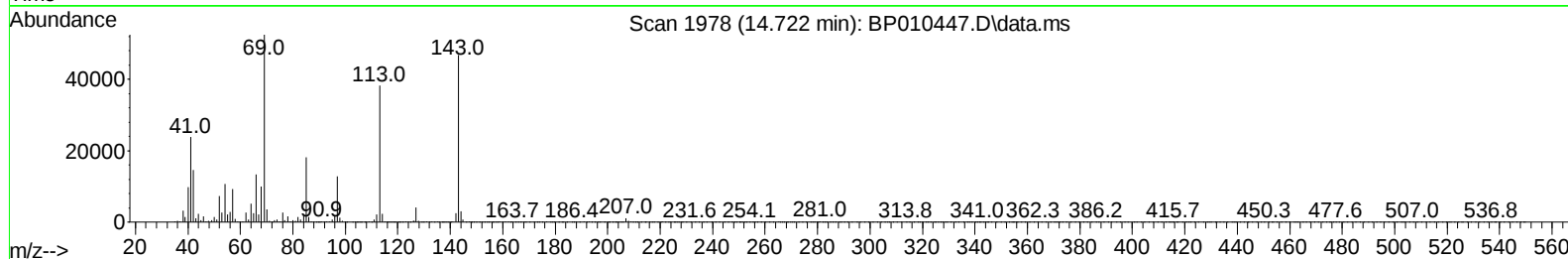
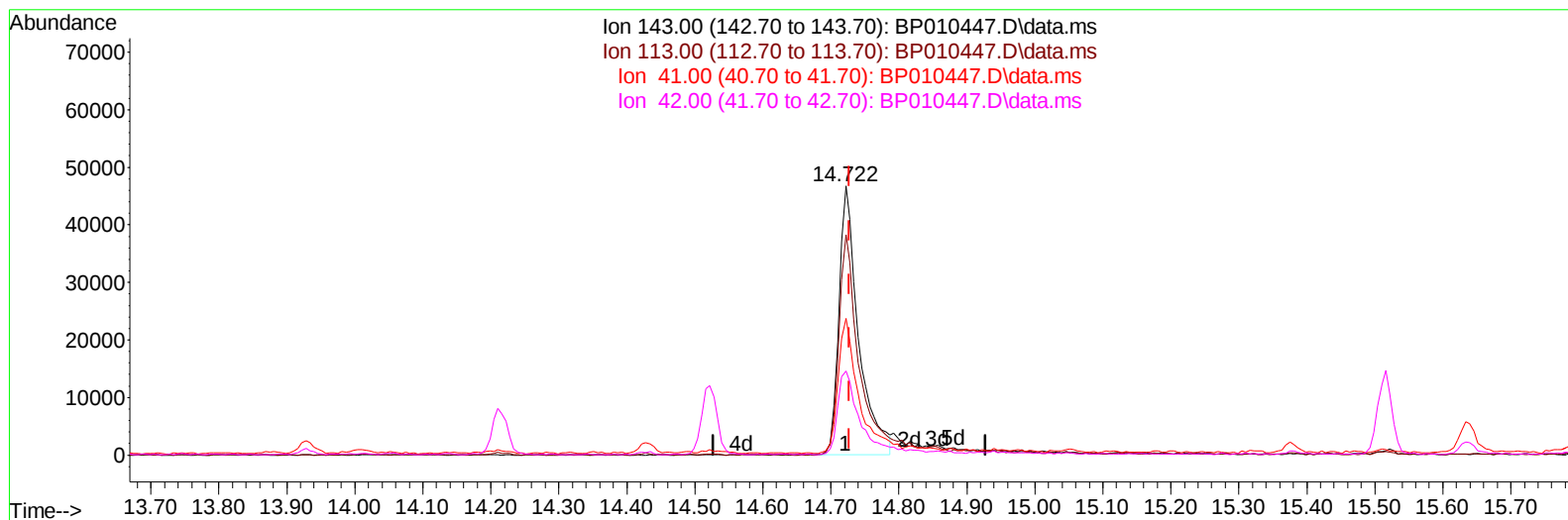
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TIC: BP010447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 13.57 ng/u1

response 93437

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.83#
41.00	23.20	50.92#
42.00	18.00	31.21#

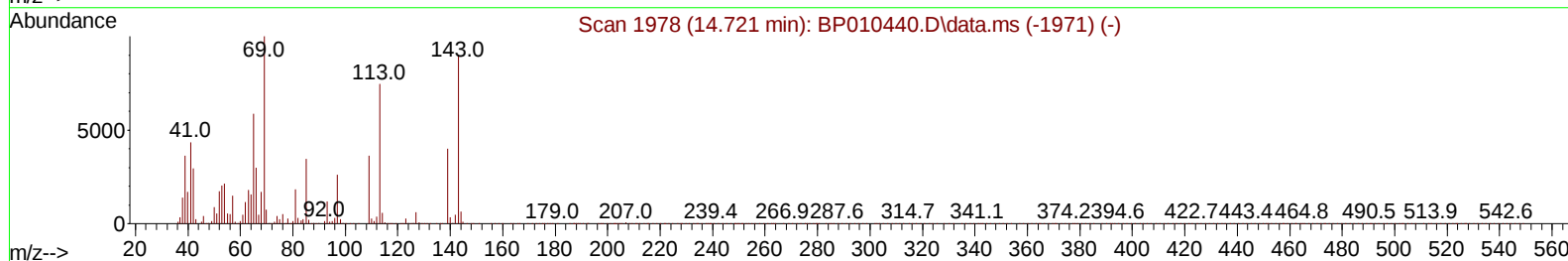
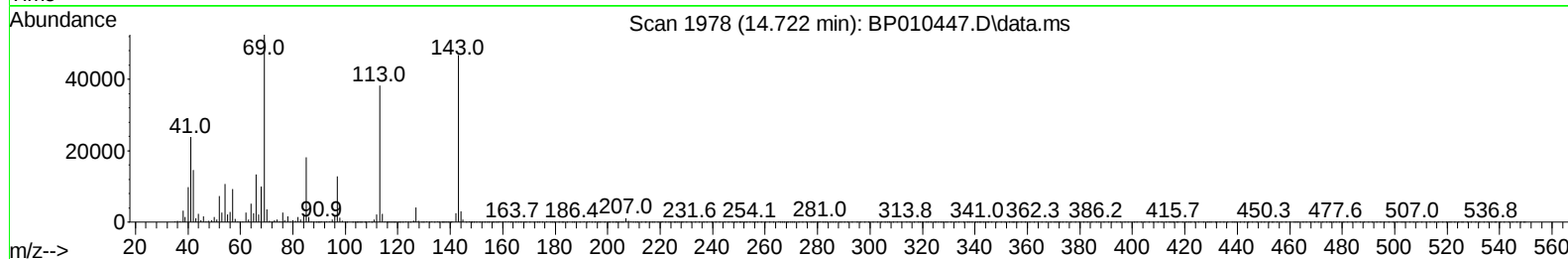
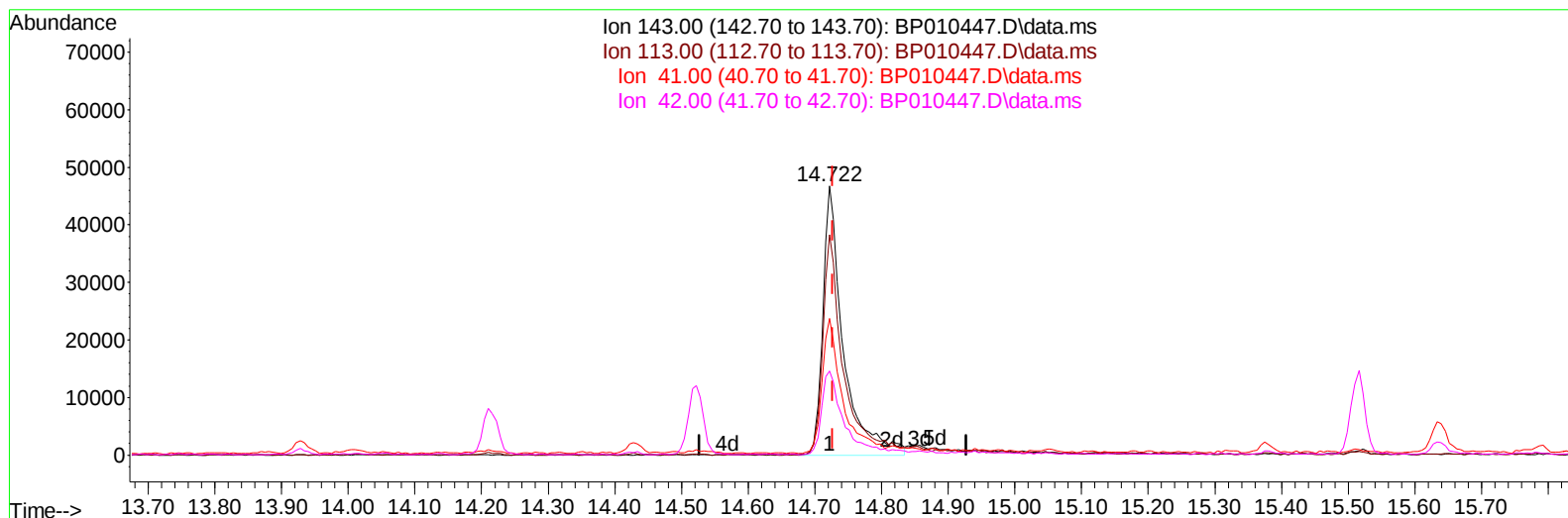
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Instrument :
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TIC: BP010447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 14.53 ng/ul m

response 100052

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.83#
41.00	23.20	50.92#
42.00	18.00	31.21#

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 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	168355	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	760277	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.522	164	525000	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1129318	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	1090680	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	1012112	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	12597m	3.189	ng/uL	0.00
4) Pyridine-d5	3.770	84	54378	4.849	ng/ul	0.00
7) Phenol-d5	7.058	99	258379	18.285	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	177735	18.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	209529	19.378	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	217759	18.134	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	110592	18.795	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	119310	19.133	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	216993	18.404	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	295309	16.845	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	800198	20.687	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	899069	20.151	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	100052m	14.532	ng/ul	0.00
60) Fluorene-d10	15.516	176	703763	20.913	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	89588	13.564	ng/ul	0.00
73) Anthracene-d10	17.375	188	1069698	20.918	ng/ul	0.00
81) Pyrene-d10	19.604	212	1419599	24.509	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.622	264	1192191	23.458	ng/ul	-0.01

Target Compounds Qvalue

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

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