

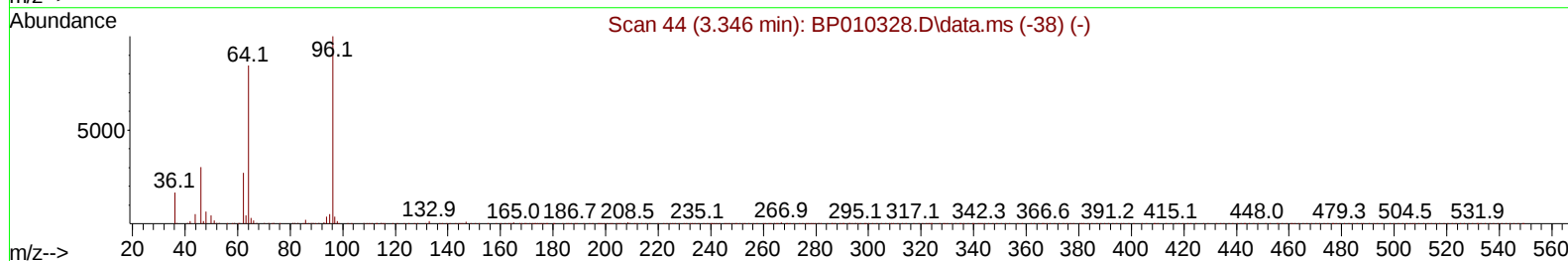
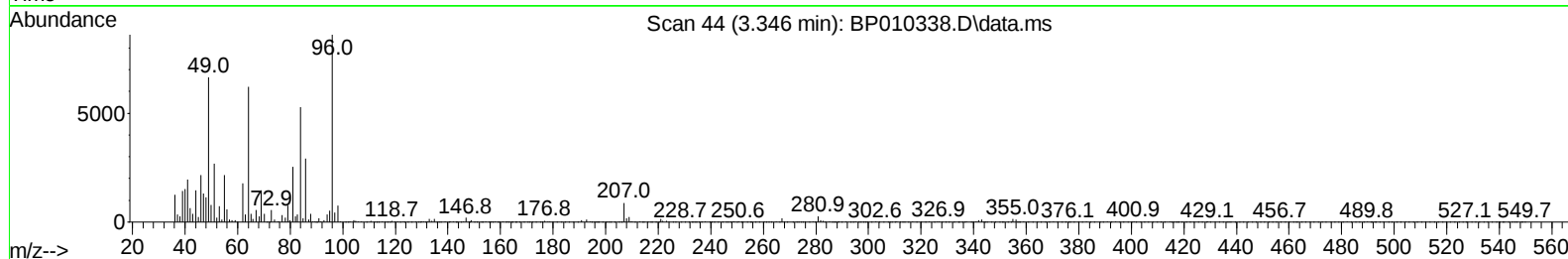
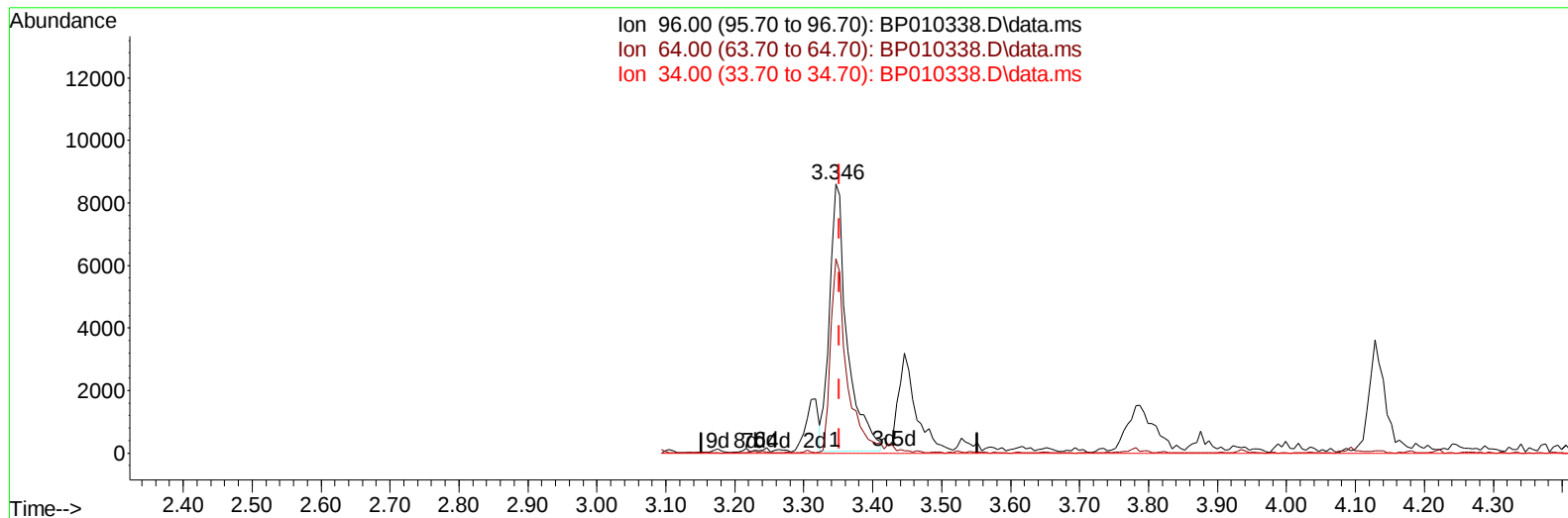
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXLB5

Manual IntegrationsAPPROVED

Quant Time: May 13 01:56:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/16/2022
Supervised By :Yogesh Patel 05/17/2022



TIC: BP010338.D\data.ms

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

3.346min (-0.006) 5.58 ng/uL

response 15321

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.25#
34.00	0.00	0.00
0.00	0.00	0.00

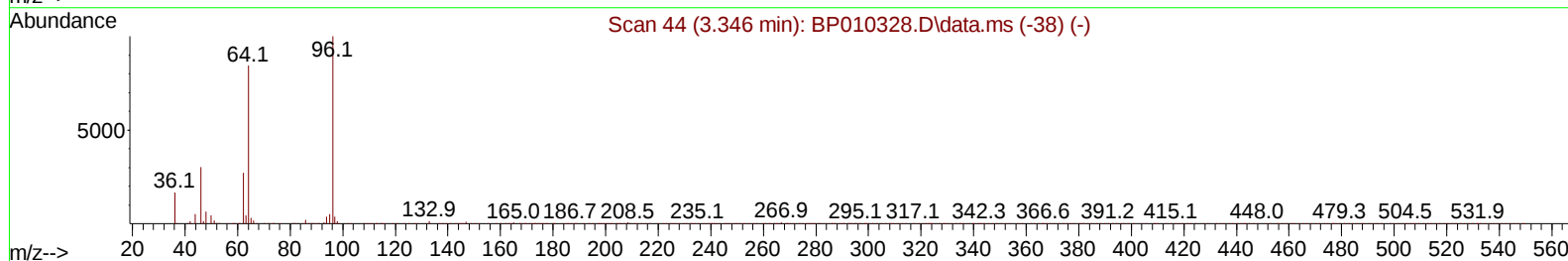
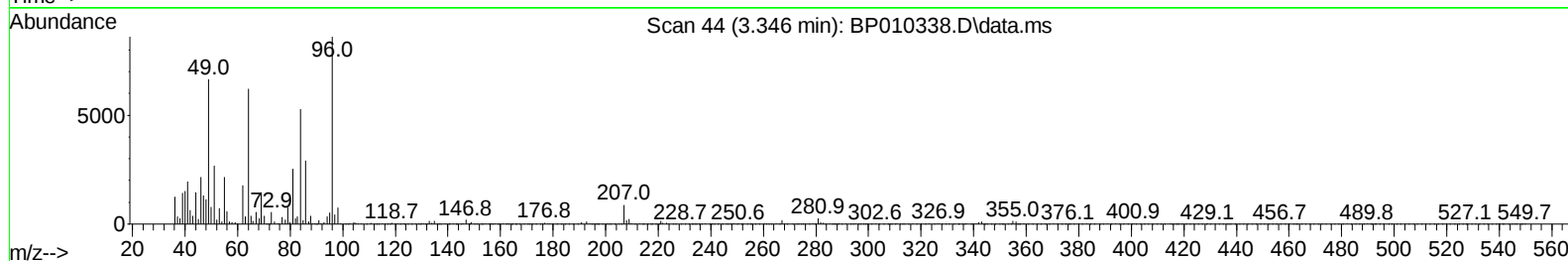
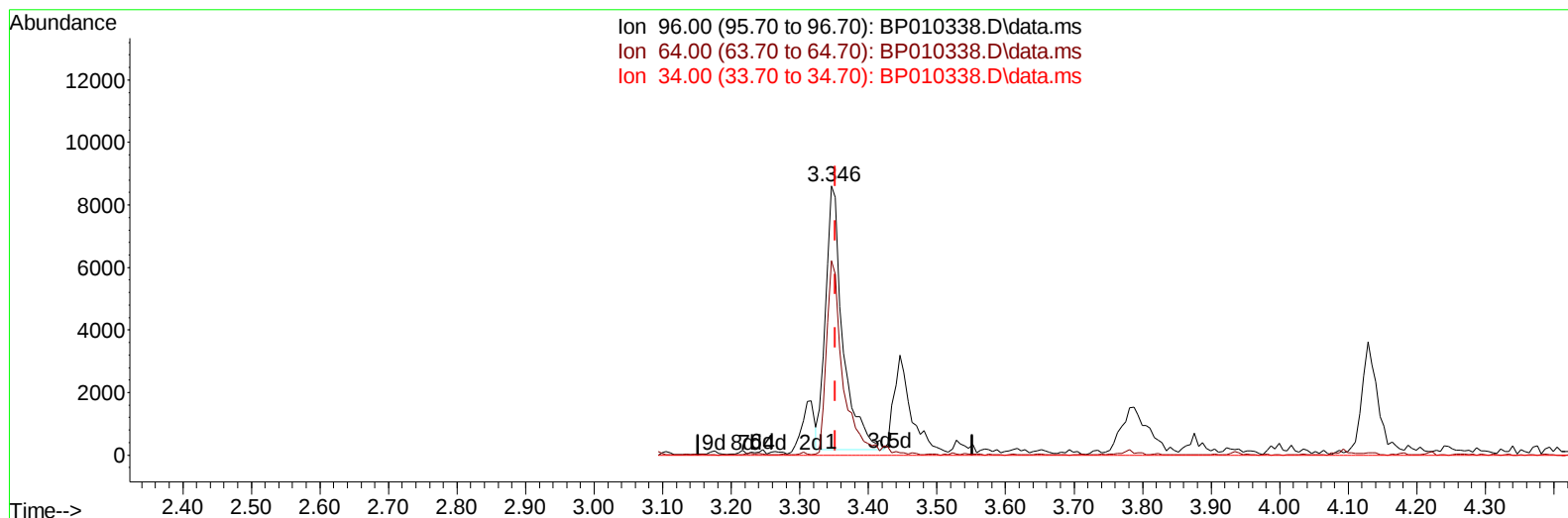
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TIC: BP010338.D\data.ms

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

3.346min (-0.006) 5.40 ng/uL m

response 14820

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.25#
34.00	0.00	0.00
0.00	0.00	0.00

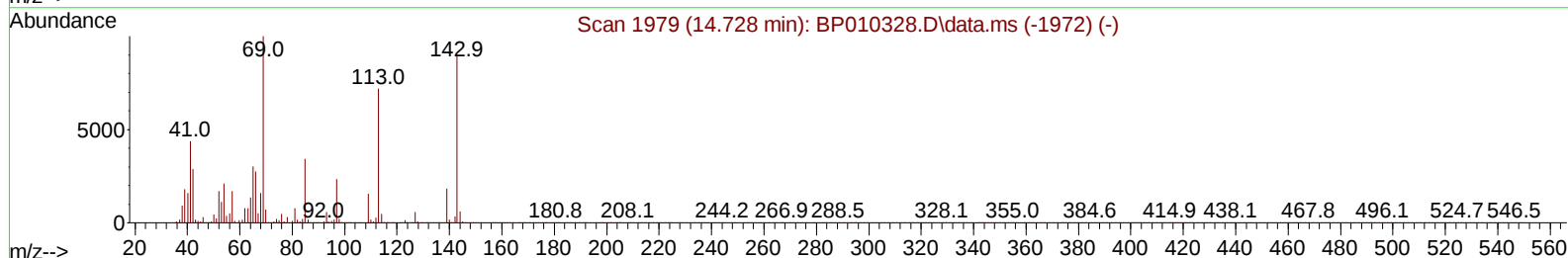
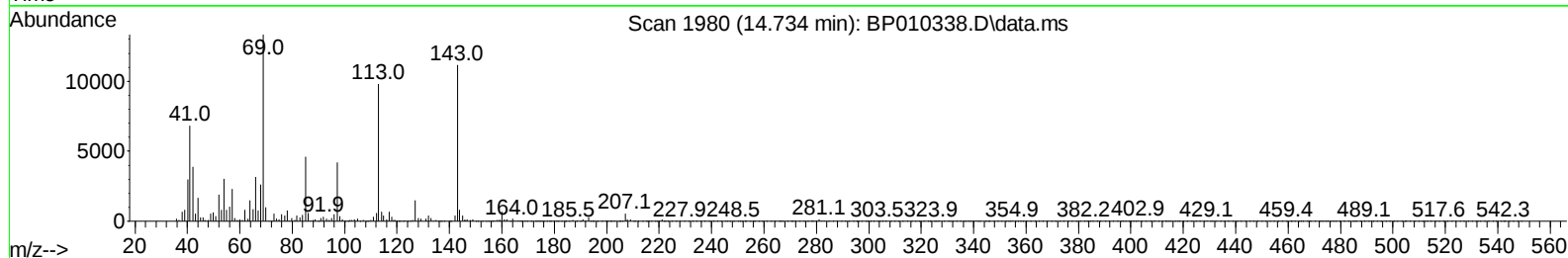
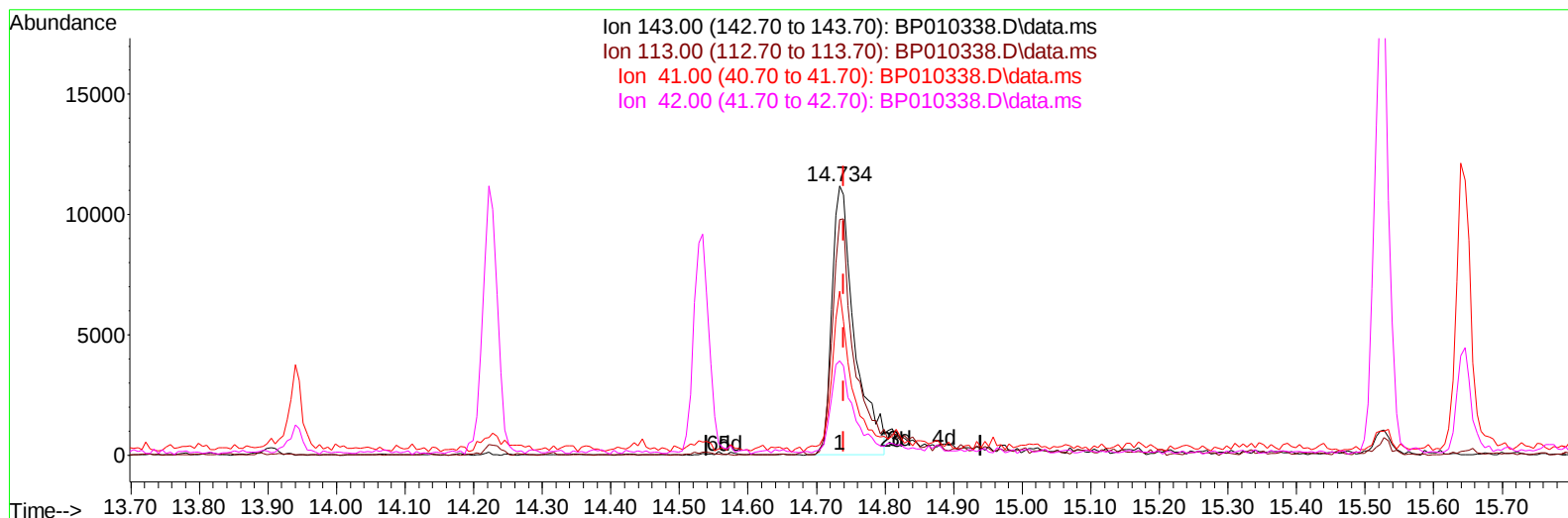
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TIC: BP010338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 4.47 ng/u1

response 25656

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.58#
41.00	23.20	61.06#
42.00	18.00	34.95#

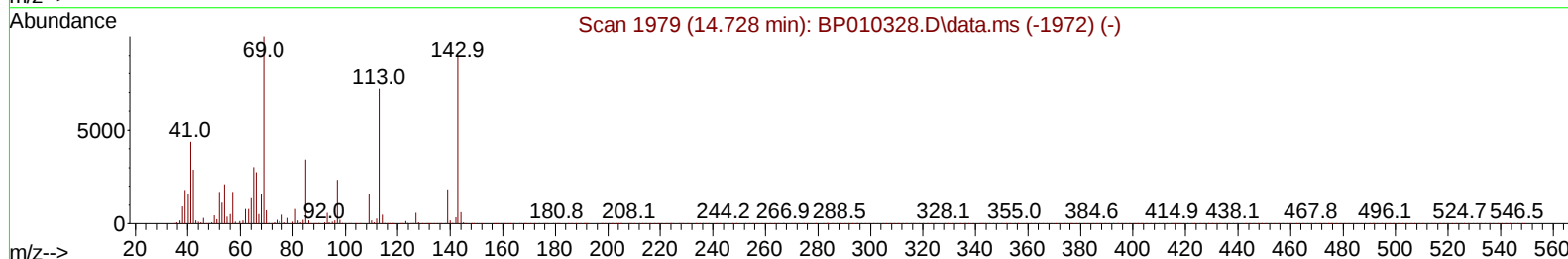
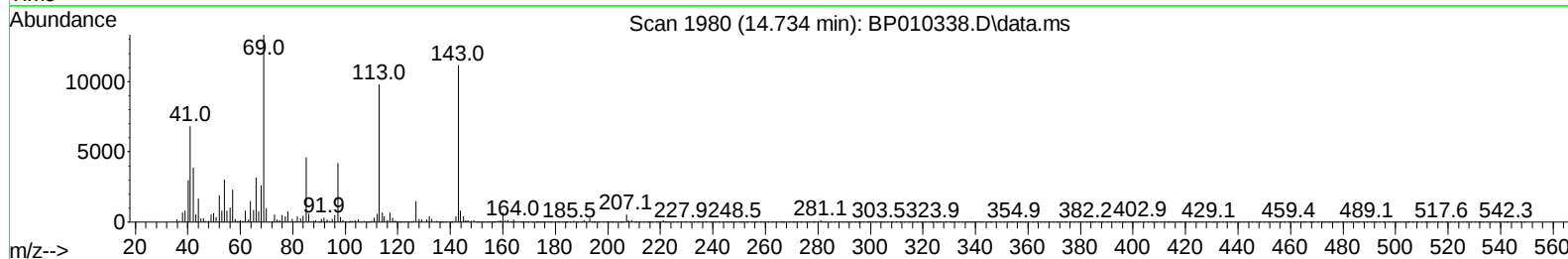
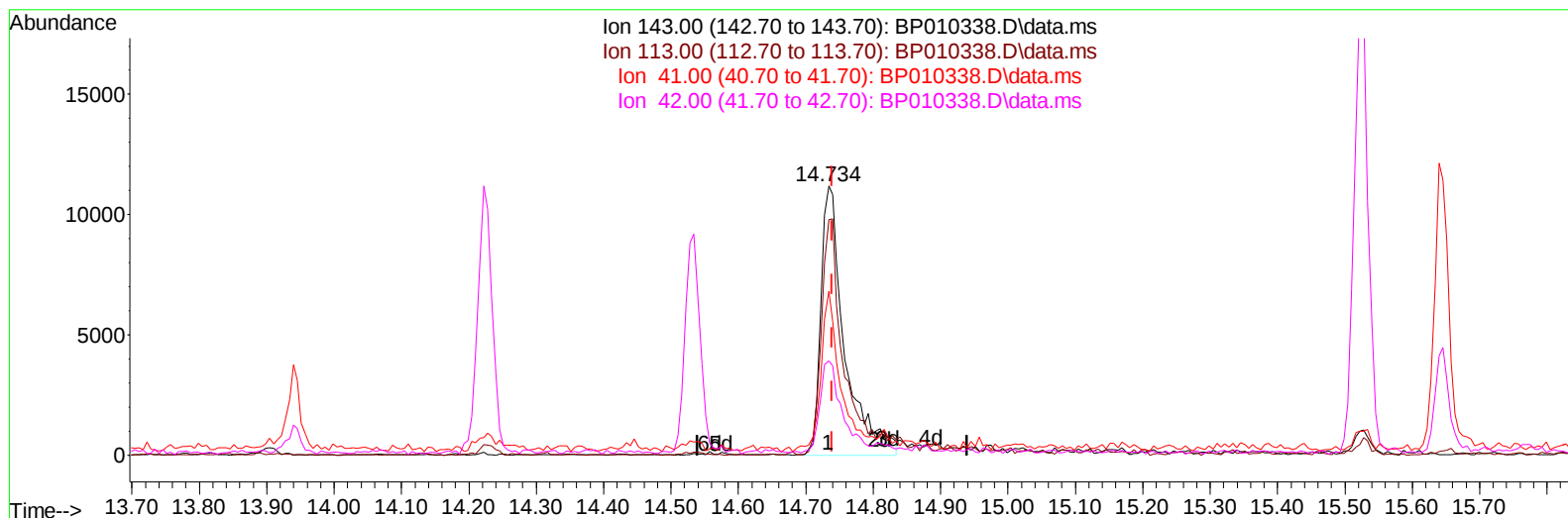
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 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP010338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 4.78 ng/ul m

response 27443

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.58#
41.00	23.20	61.06#
42.00	18.00	34.95#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	116691	20.000	ng/ul	0.00
20) Naphthalene-d8	10.705	136	520926	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	375576	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	840230	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.363	240	889292	20.000	ng/ul	#-0.01
88) Perylene-d12	23.792	264	853803	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	14820m	5.395	ng/uL	0.00
4) Pyridine-d5	3.776	84	58975	7.715	ng/ul	0.00
7) Phenol-d5	7.069	99	60239	5.558	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	230039	34.310	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	196221	24.583	ng/ul	0.00
15) 4-Methylphenol-d8	8.611	113	129516	14.535	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	147436	35.298	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	146500	30.762	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	244677	28.228	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	347321	27.142	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1080188	36.101	ng/ul	-0.01
49) Acenaphthylene-d8	14.228	160	1155220	32.383	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	27443m	4.782	ng/ul	0.00
60) Fluorene-d10	15.528	176	931813	36.894	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.646	200	173905	30.224	ng/ul	-0.01
73) Anthracene-d10	17.381	188	1612689	39.194	ng/ul	-0.01
81) Pyrene-d10	19.616	212	2013913	40.981	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1880276	41.294	ng/ul	-0.02

Target Compounds Qvalue

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

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