

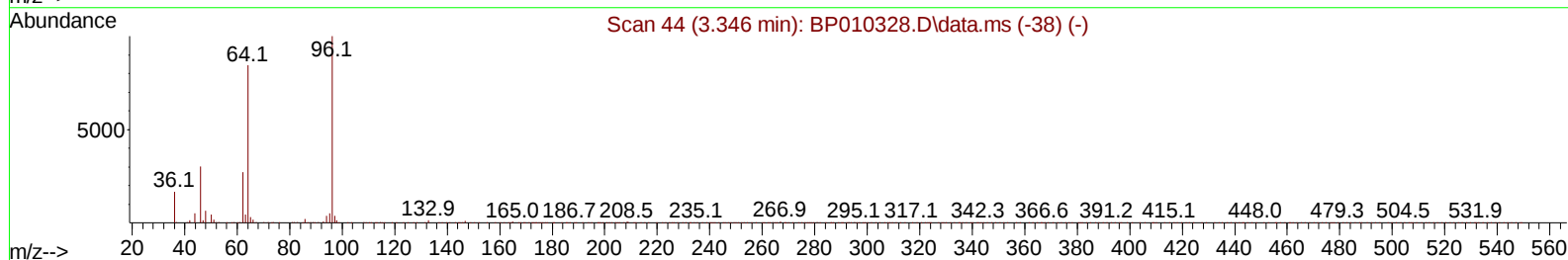
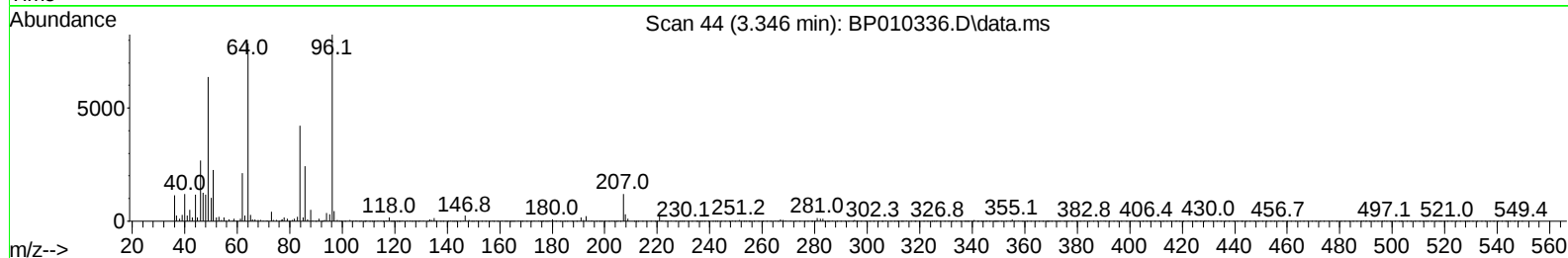
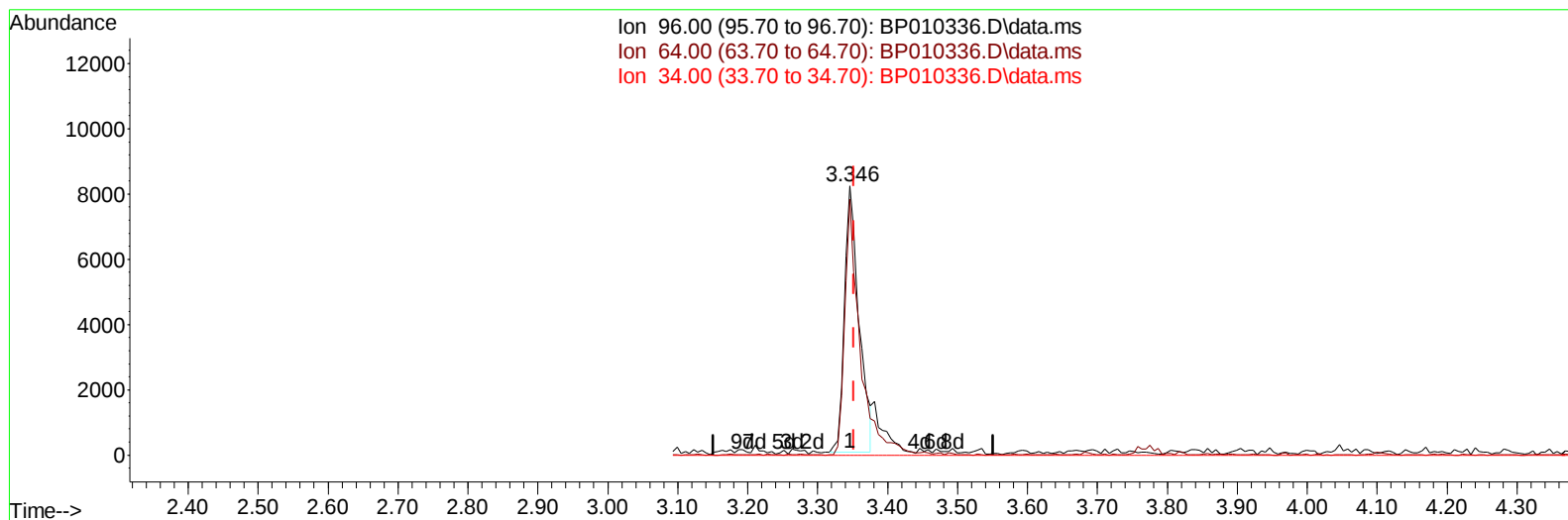
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010336.D
 Acq On : 12 May 2022 22:31
 Operator : CG/JU
 Sample : N2746-05
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBSL1

Manual IntegrationsAPPROVED

Quant Time: May 13 00:04:07 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: BP010336.D\data.ms

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

3.346min (-0.006) 4.41 ng/uL

response 12007

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	95.10#
34.00	0.00	0.00
0.00	0.00	0.00

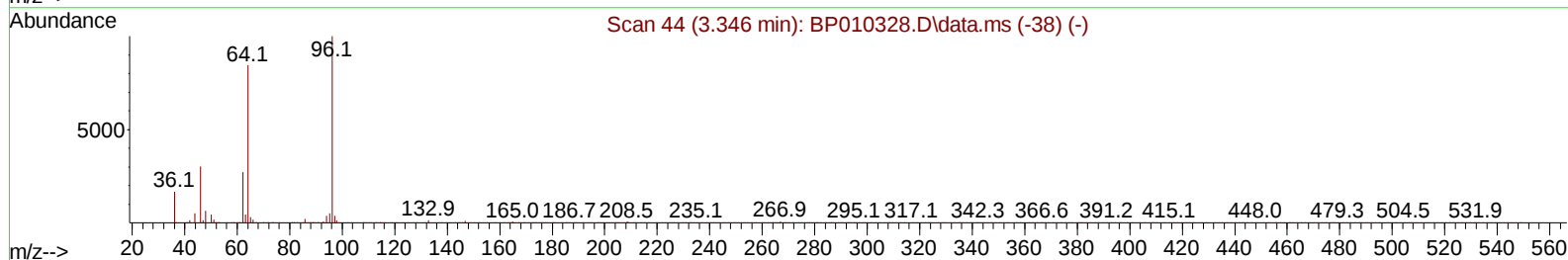
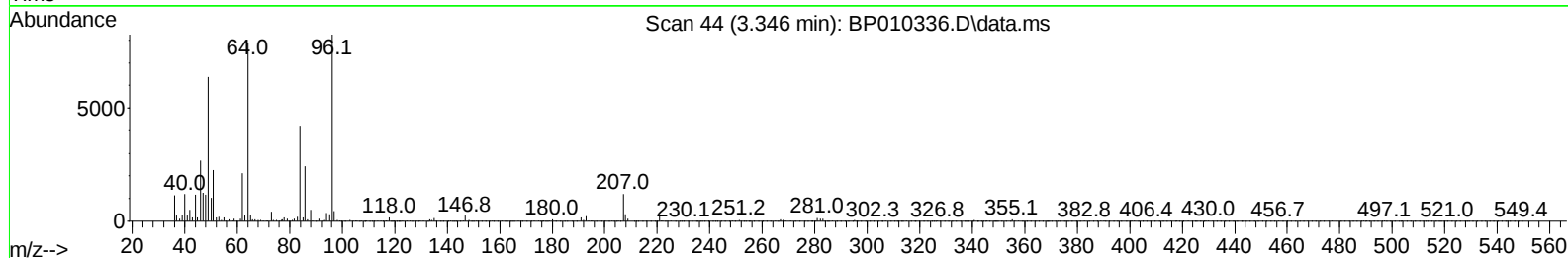
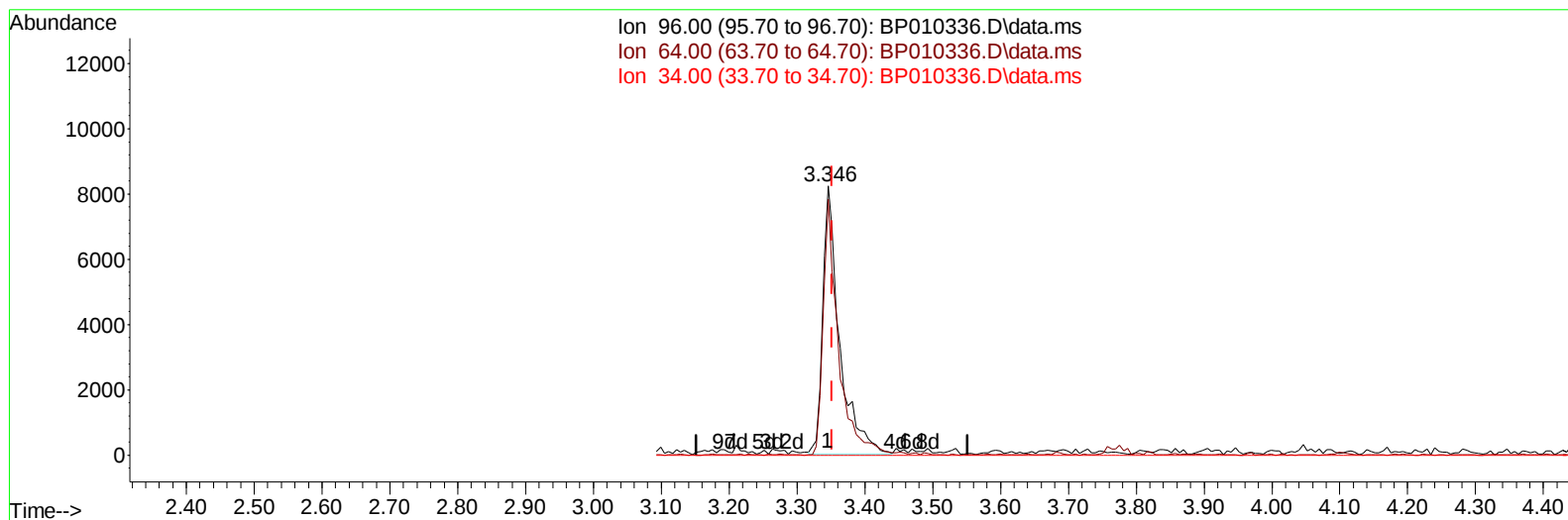
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TIC: BP010336.D\data.ms

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

3.346min (-0.006) 5.14 ng/uL m

response 14001

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	95.10#
34.00	0.00	0.00
0.00	0.00	0.00

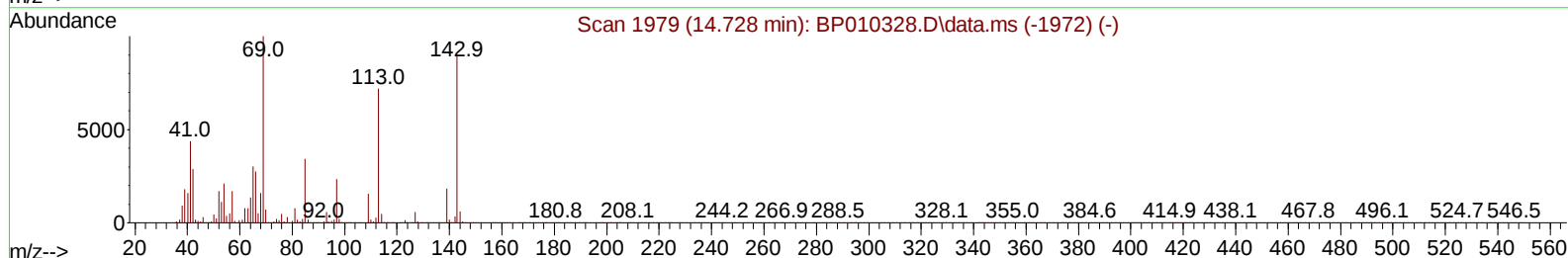
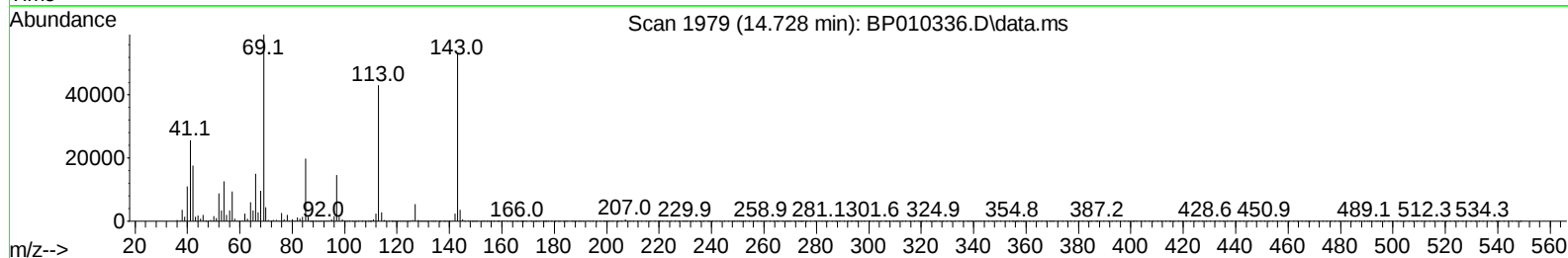
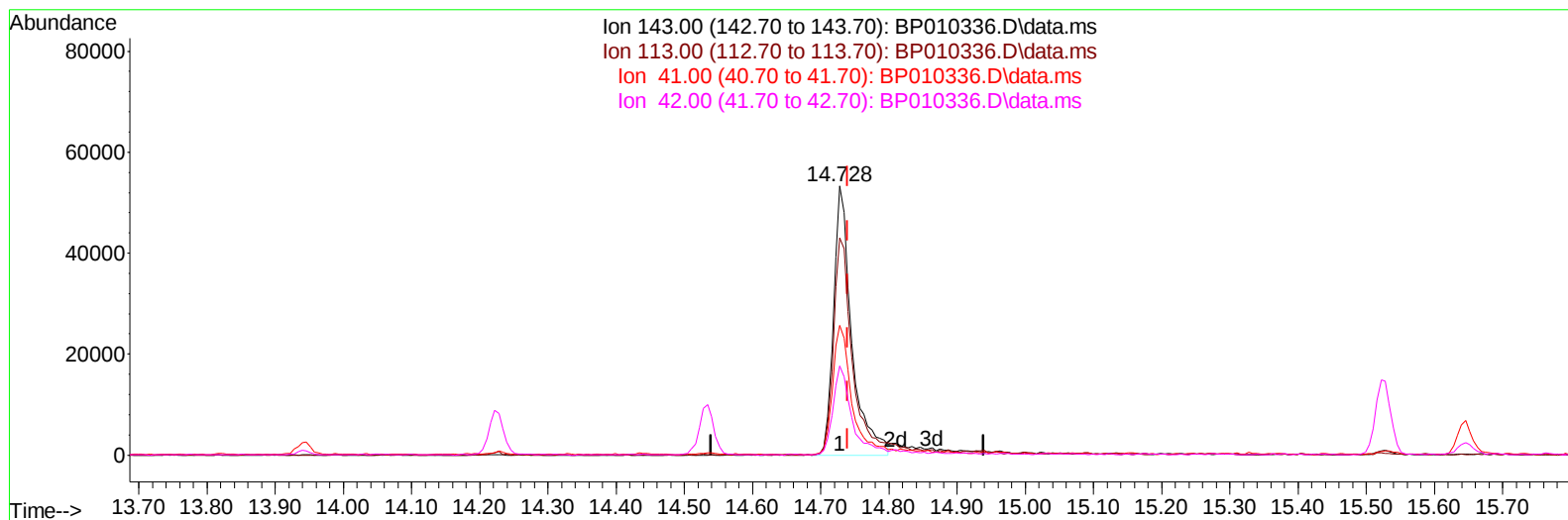
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TIC: BP010336.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.012) 16.82 ng/u1

response 98960

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.81#
41.00	23.20	48.16#
42.00	18.00	33.11#

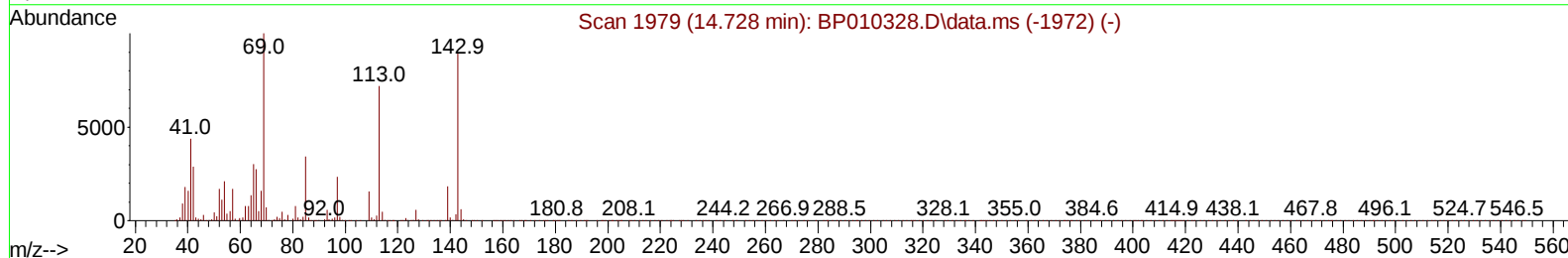
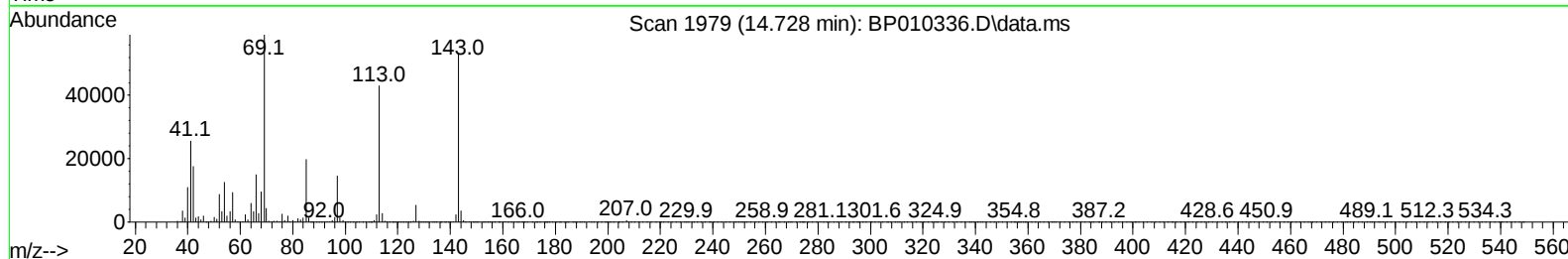
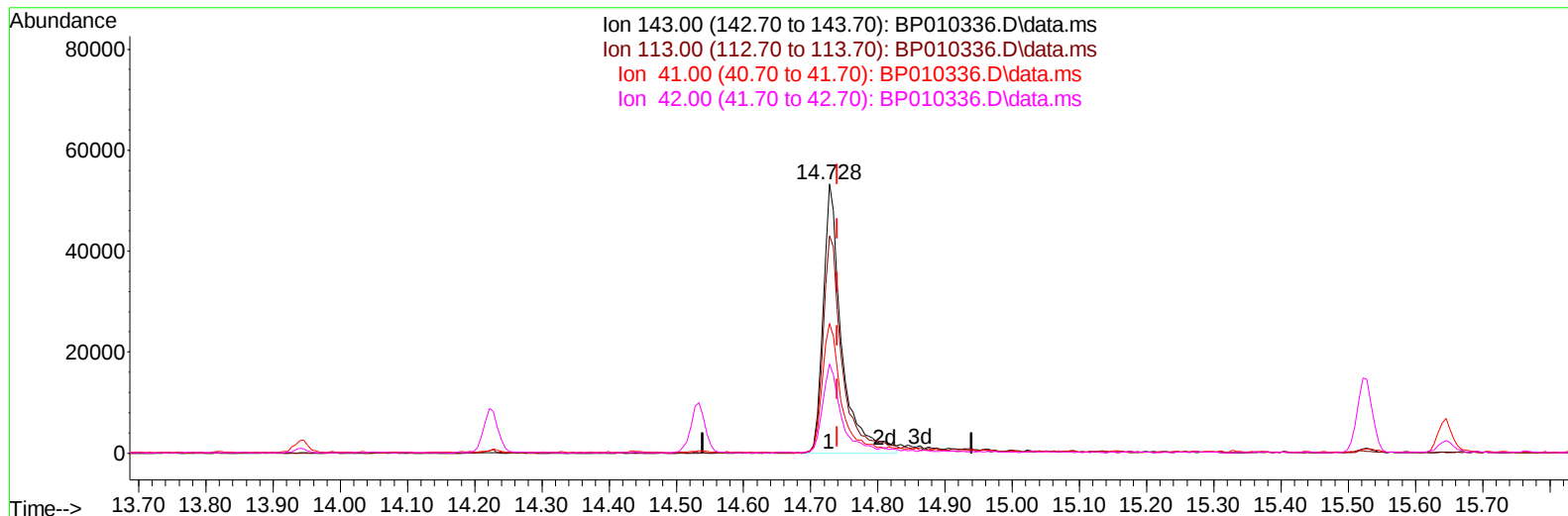
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Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010336.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.012) 17.41 ng/ul m

response 102458

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.81#
41.00	23.20	48.16#
42.00	18.00	33.11#

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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	115721	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	529554	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	385074	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	863324	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.369	240	917939	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	907822	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	14001m	5.140	ng/uL	0.00
4) Pyridine-d5	3.764	84	200716	26.479	ng/ul	0.00
7) Phenol-d5	7.069	99	272690	25.371	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	190950	28.719	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	217037	27.419	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	237498	26.877	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	118444	27.895	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	127703	26.378	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	228959	25.984	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.840	131	336959	25.903	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	868434	28.308	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	962768	26.323	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	102458m	17.412	ng/ul	-0.01
60) Fluorene-d10	15.528	176	742453	28.671	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	91877	15.541	ng/ul	-0.01
73) Anthracene-d10	17.386	188	1188308	28.108	ng/ul	0.00
81) Pyrene-d10	19.616	212	1479311	29.163	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1401358	28.945	ng/ul	-0.02

Target Compounds Qvalue

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

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