

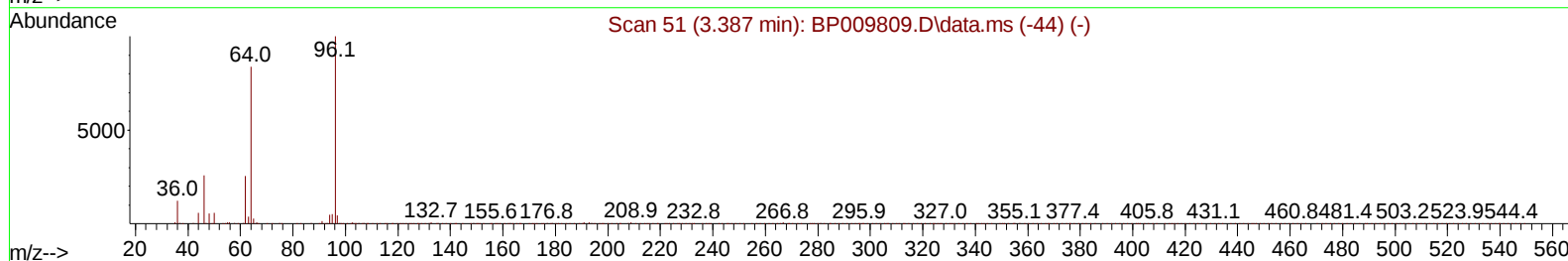
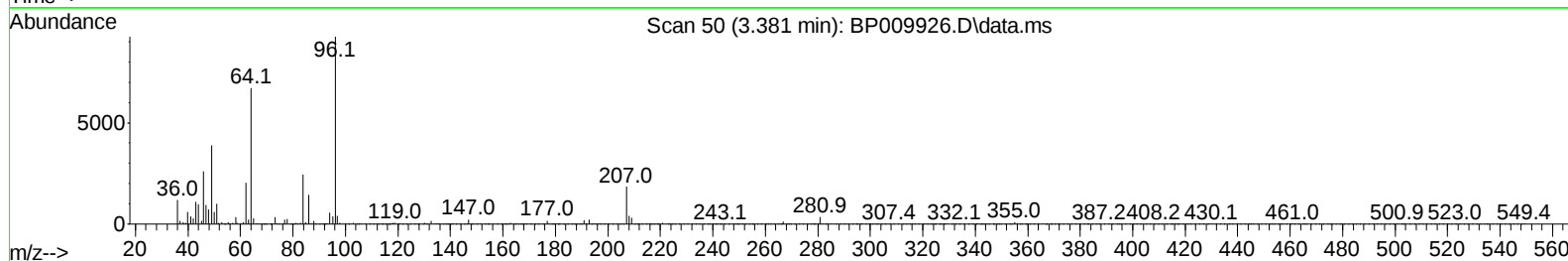
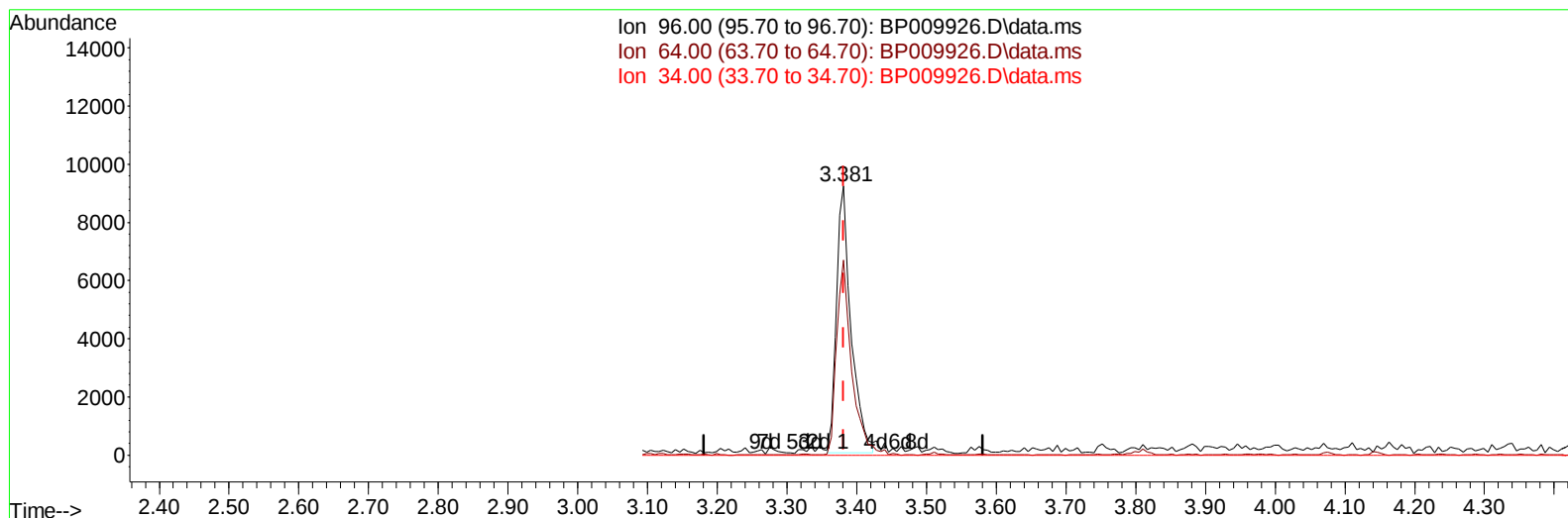
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009926.D
 Acq On : 18 Apr 2022 21:11
 Operator : CG/JU
 Sample : N2421-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY8

Manual IntegrationsAPPROVED

Quant Time: Apr 19 00:58:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009926.D\data.ms

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

3.381min (-0.000) 4.88 ng/uL

response 13263

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.51#
34.00	0.00	0.00
0.00	0.00	0.00

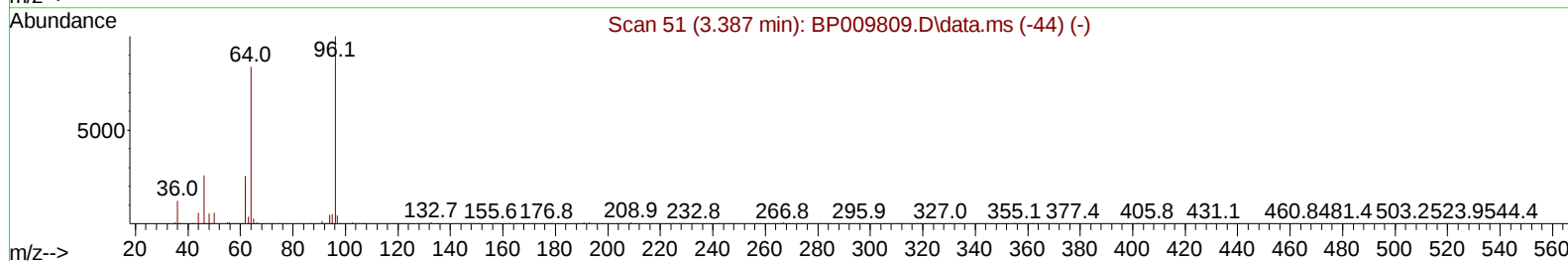
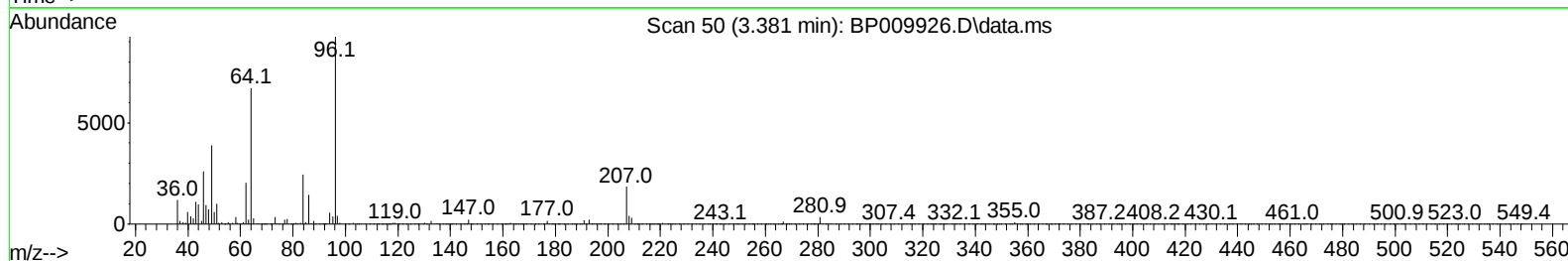
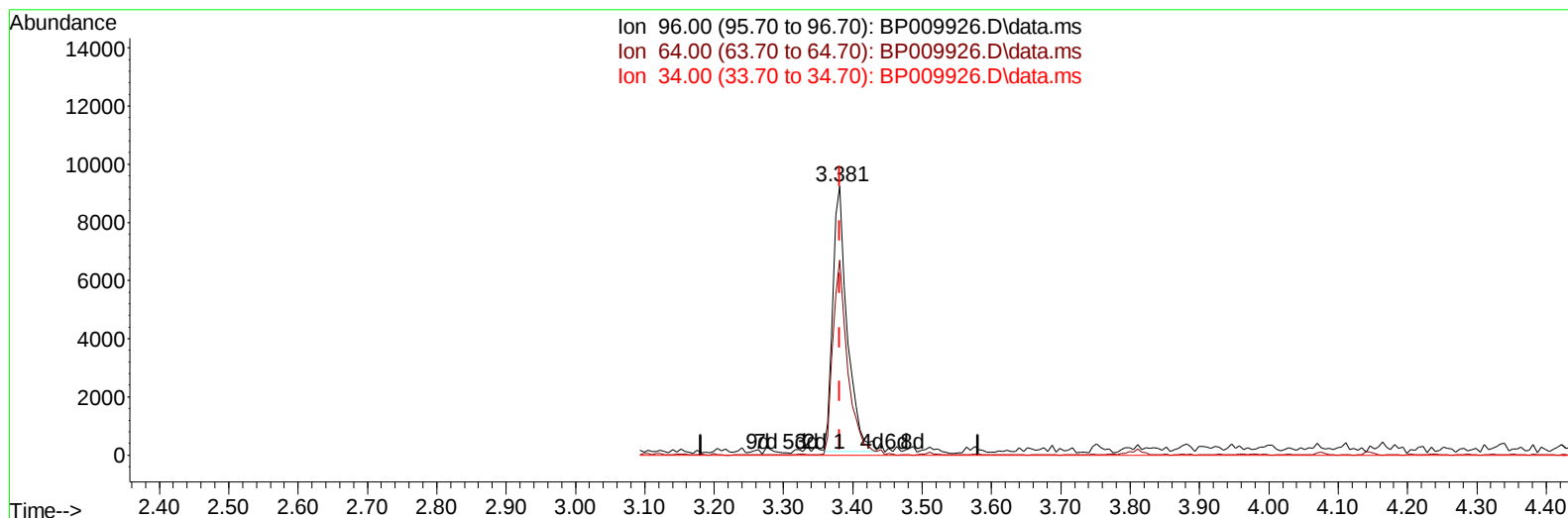
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TIC: BP009926.D\data.ms

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

3.381min (-0.000) 4.86 ng/uL m

response 13214

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.51#
34.00	0.00	0.00
0.00	0.00	0.00

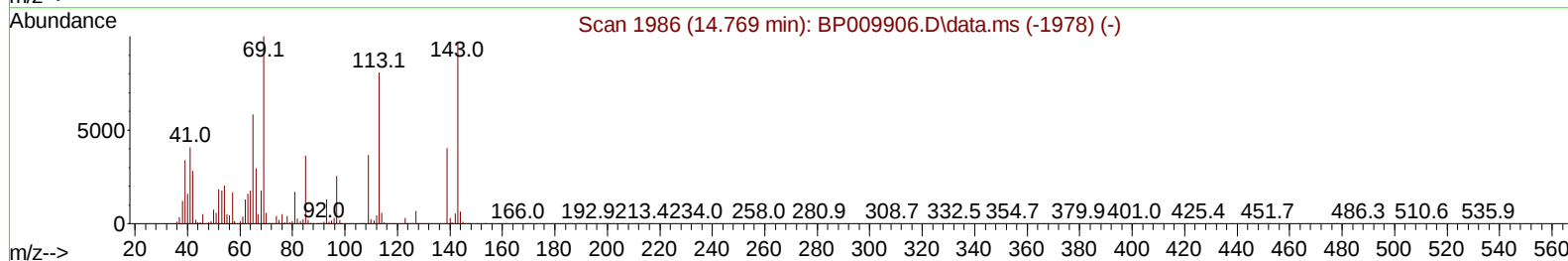
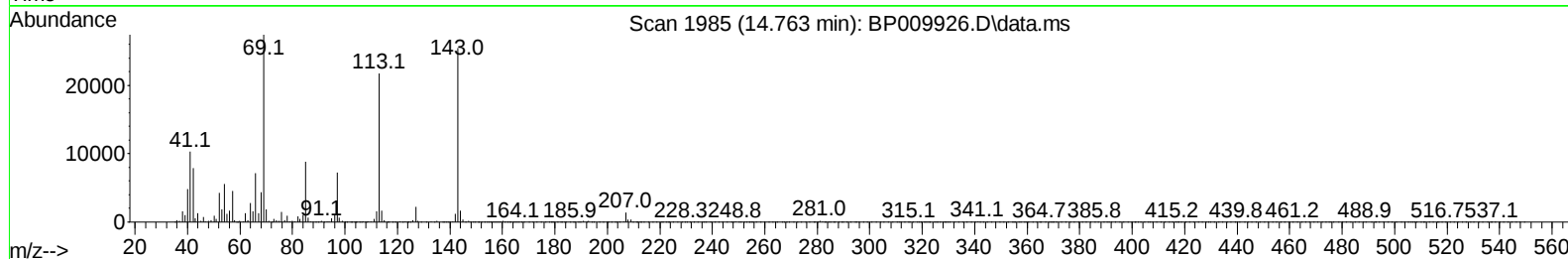
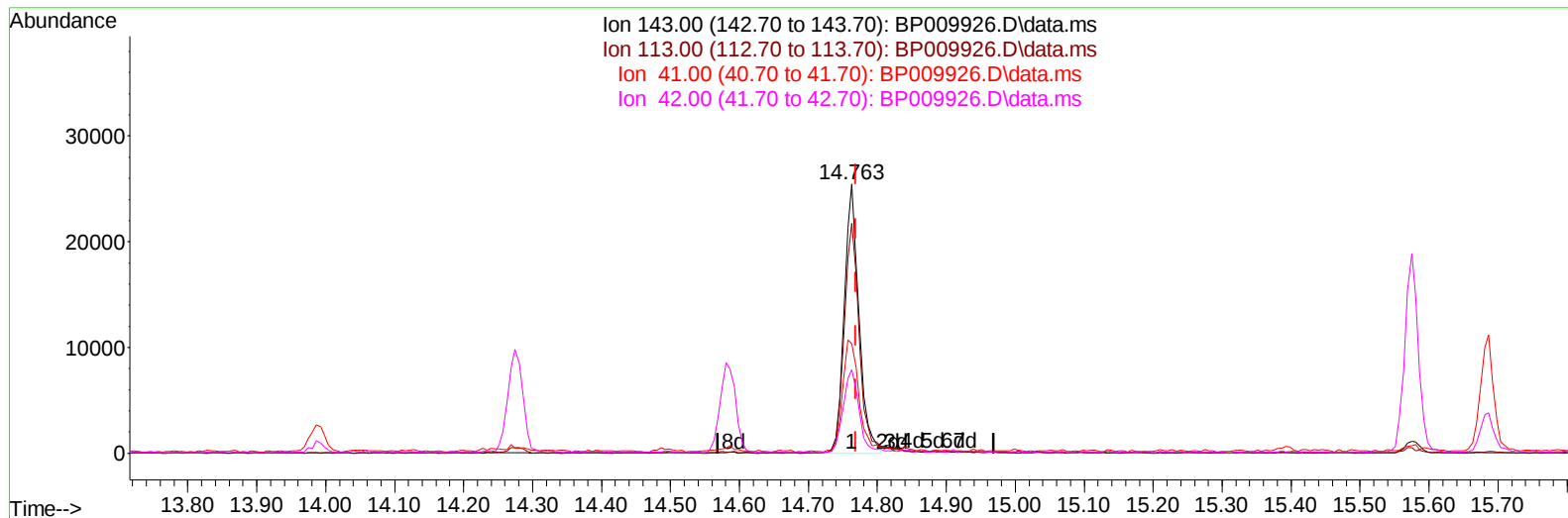
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TIC: BP009926.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 6.56 ng/u1

response 37473

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.54#
41.00	23.20	40.63#
42.00	18.00	31.09#

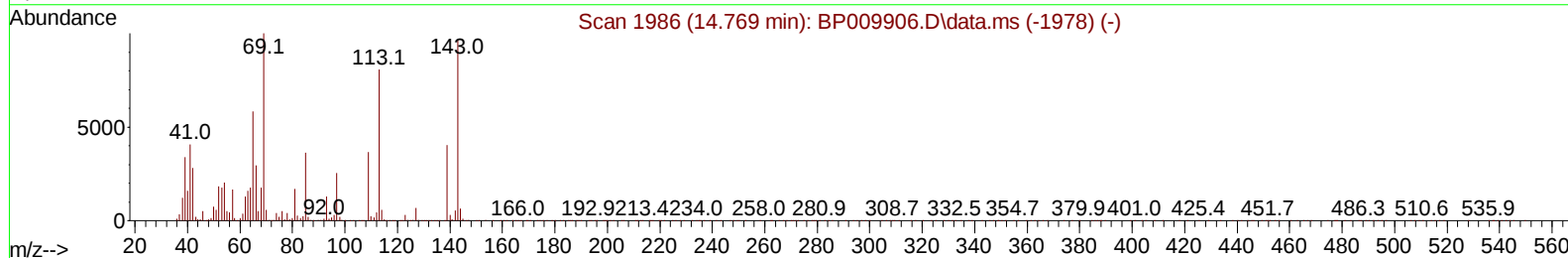
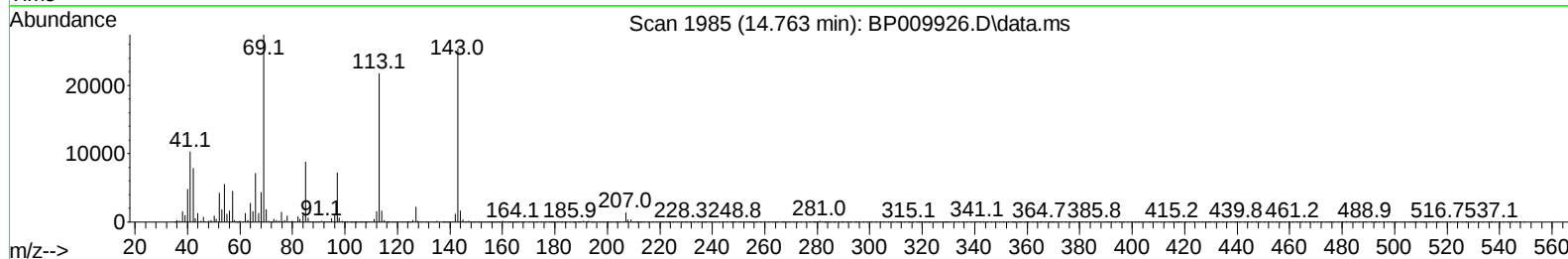
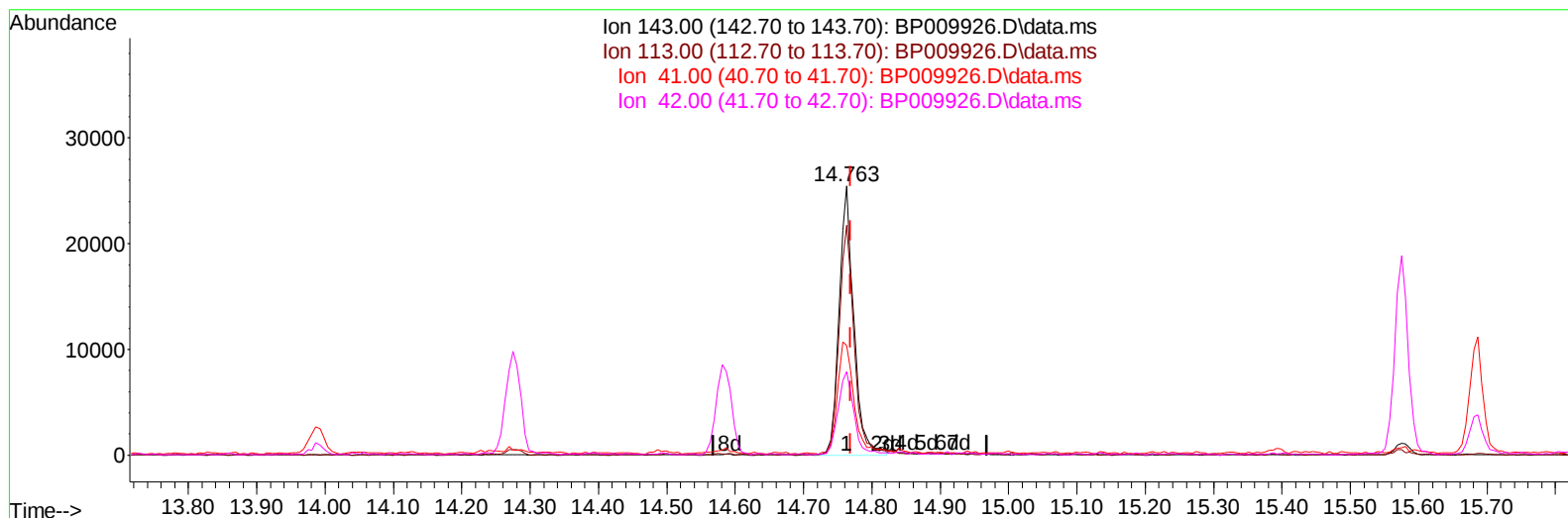
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 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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TIC: BP009926.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 6.68 ng/ul m

response 38120

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.54#
41.00	23.20	40.63#
42.00	18.00	31.09#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	121272	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	544791	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.587	164	400875	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	904462	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	900209	20.000	ng/ul	# 0.00
88) Perylene-d12	23.874	264	839002	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13214m	4.863	ng/uL	0.00
4) Pyridine-d5	3.799	84	66355	8.857	ng/ul	0.00
7) Phenol-d5	7.105	99	84485	7.935	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	204038	32.190	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	201309	25.066	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	157733	17.747	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	136185	34.665	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	137881	32.012	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	252182	27.980	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	272093	21.196	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1079648	34.286	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1145366	30.702	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	38120m	6.675	ng/ul	0.00
60) Fluorene-d10	15.575	176	930595	35.130	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	165967	30.975	ng/ul	0.00
73) Anthracene-d10	17.433	188	1626598	37.562	ng/ul	0.00
81) Pyrene-d10	19.663	212	2039607	41.485	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	1751432	40.086	ng/ul	-0.01

Target Compounds Qvalue

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

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