

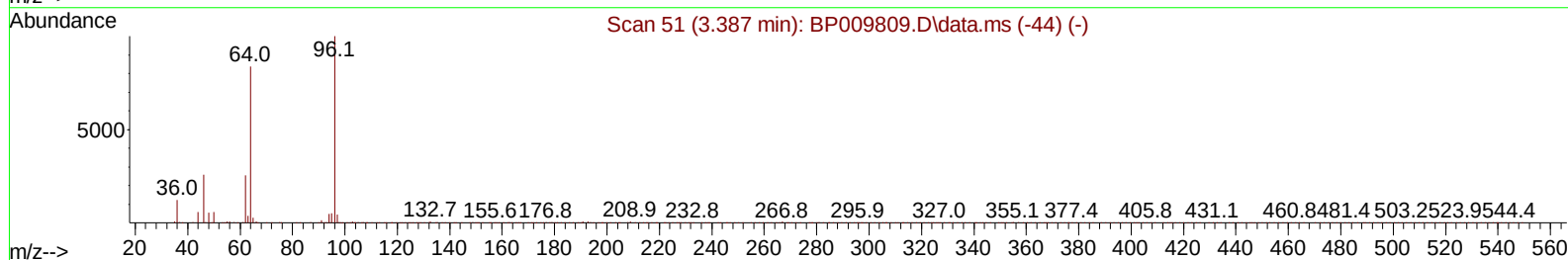
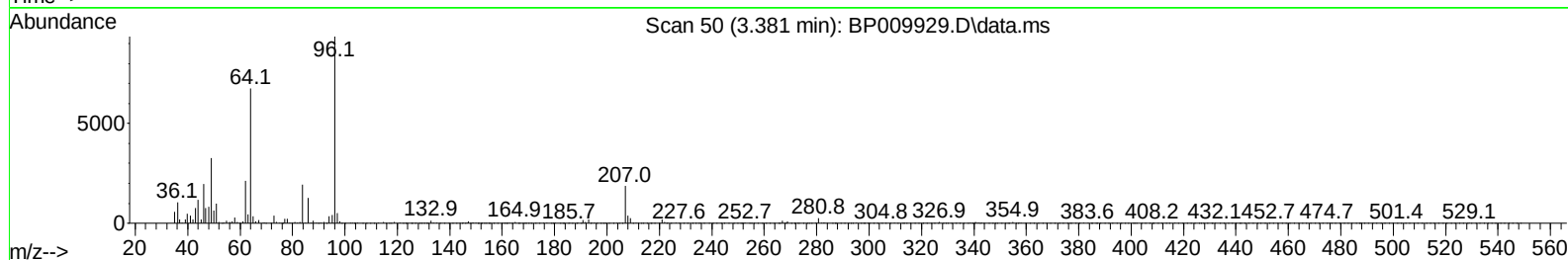
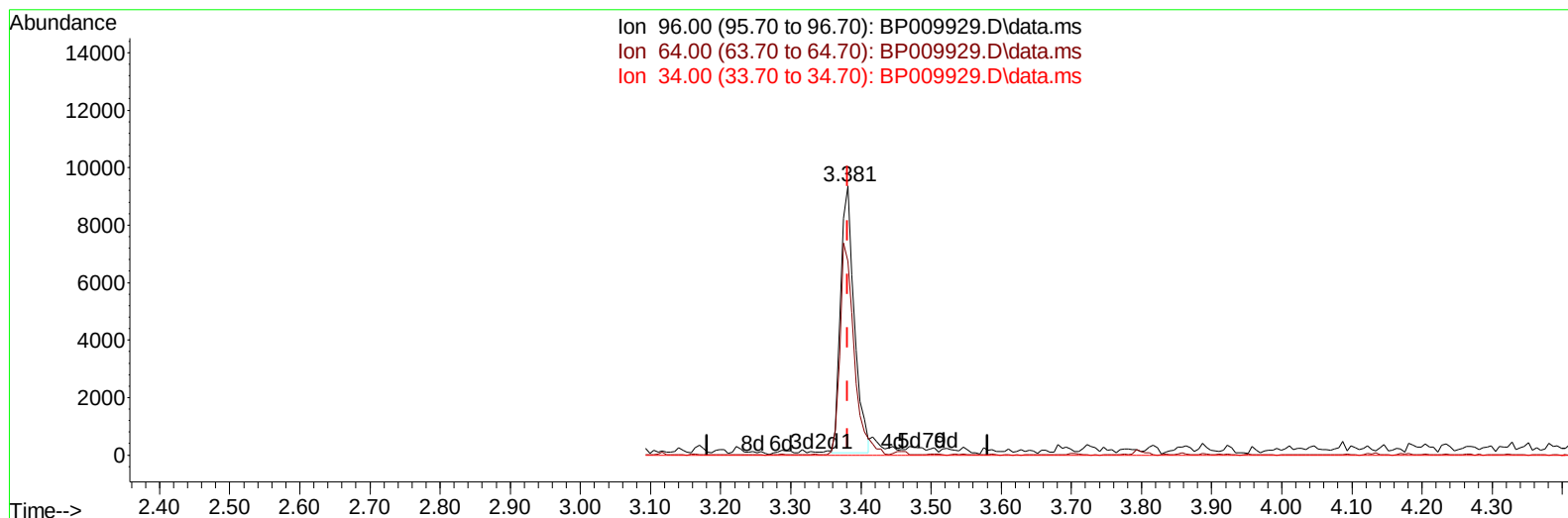
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009929.D
 Acq On : 18 Apr 2022 22:52
 Operator : CG/JU
 Sample : N2421-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J56

Manual IntegrationsAPPROVED

Quant Time: Apr 19 00:59:35 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: BP009929.D\data.ms

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

3.381min (-0.000) 4.76 ng/uL

response 12724

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.06#
34.00	0.00	0.00
0.00	0.00	0.00

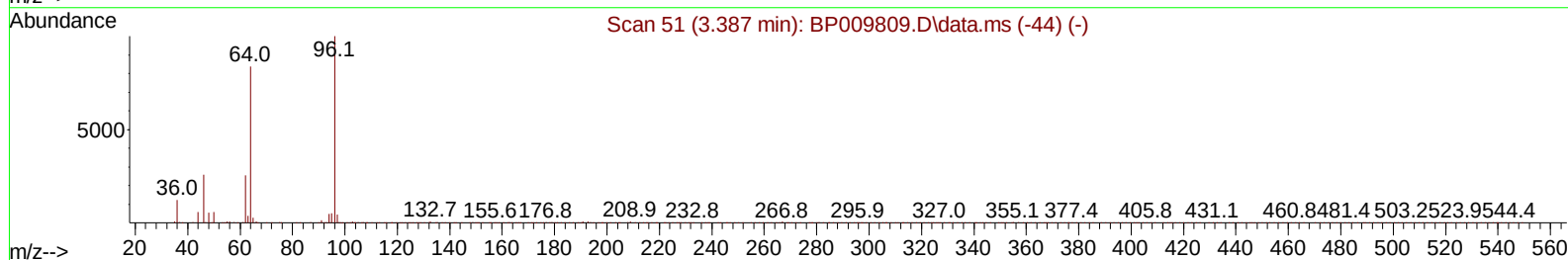
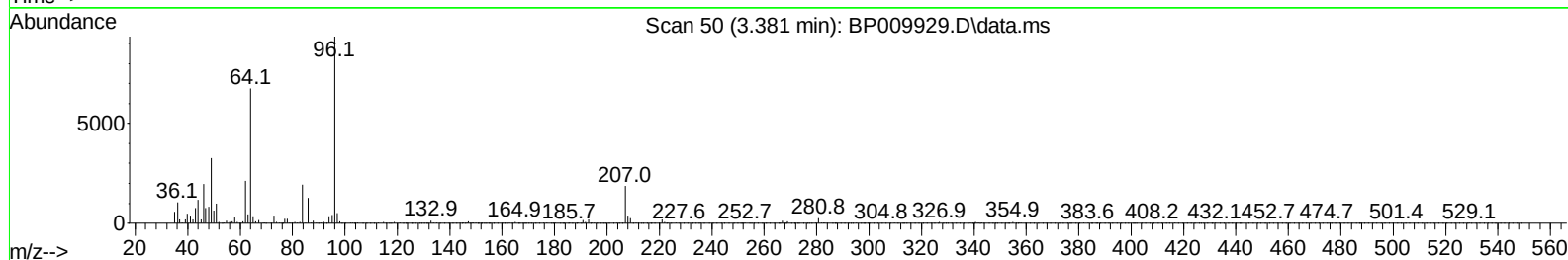
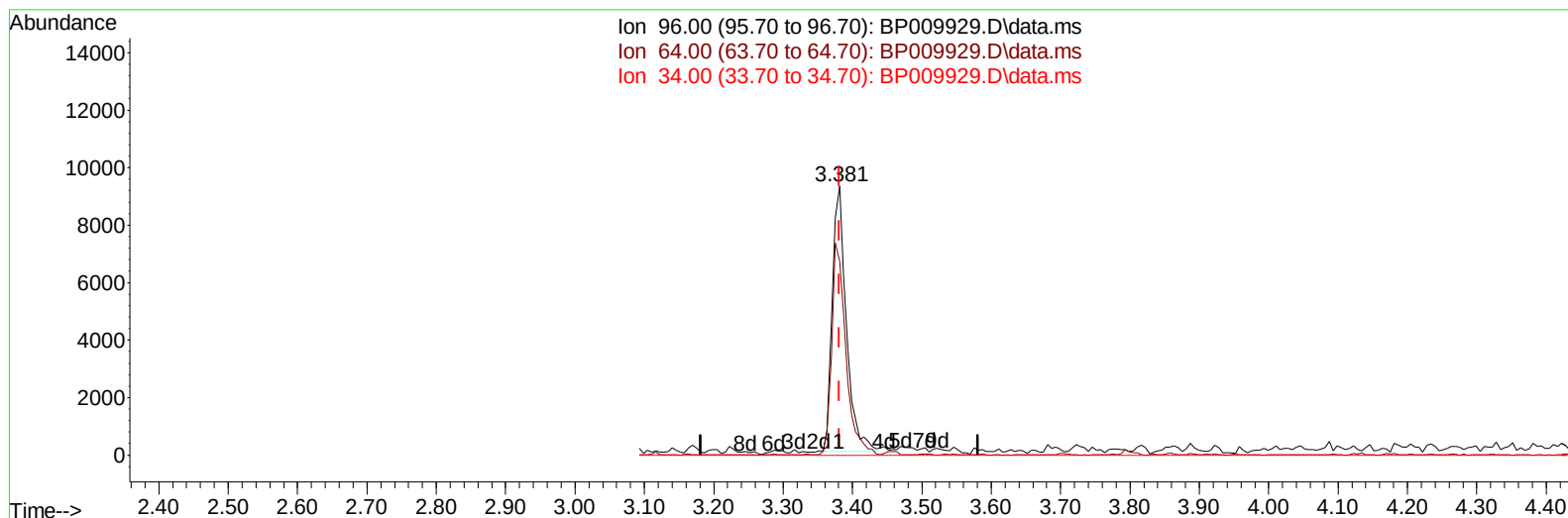
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TIC: BP009929.D\data.ms

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

3.381min (-0.000) 4.86 ng/uL m

response 12996

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.06#
34.00	0.00	0.00
0.00	0.00	0.00

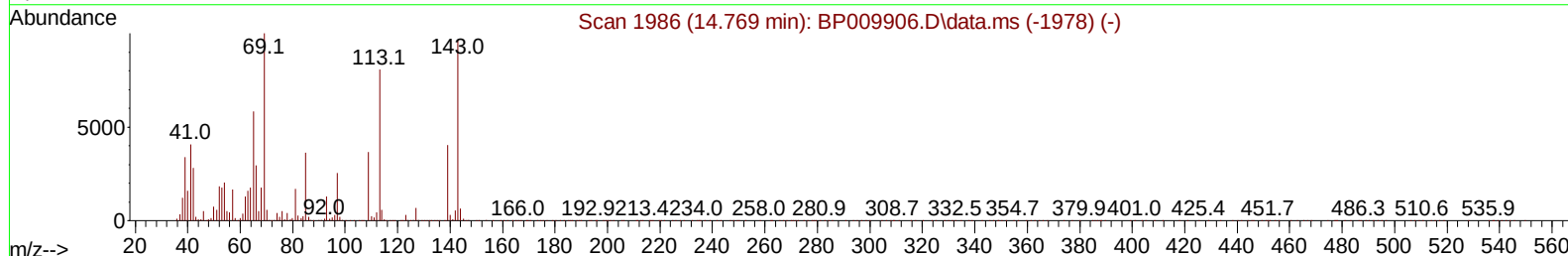
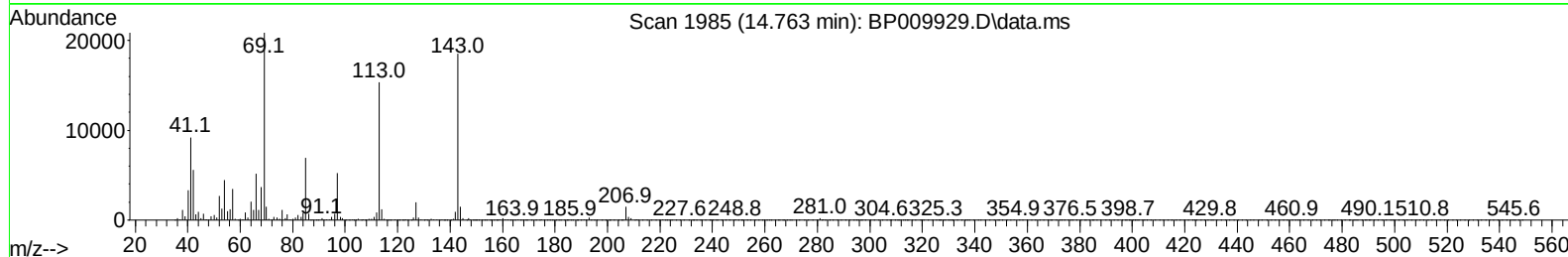
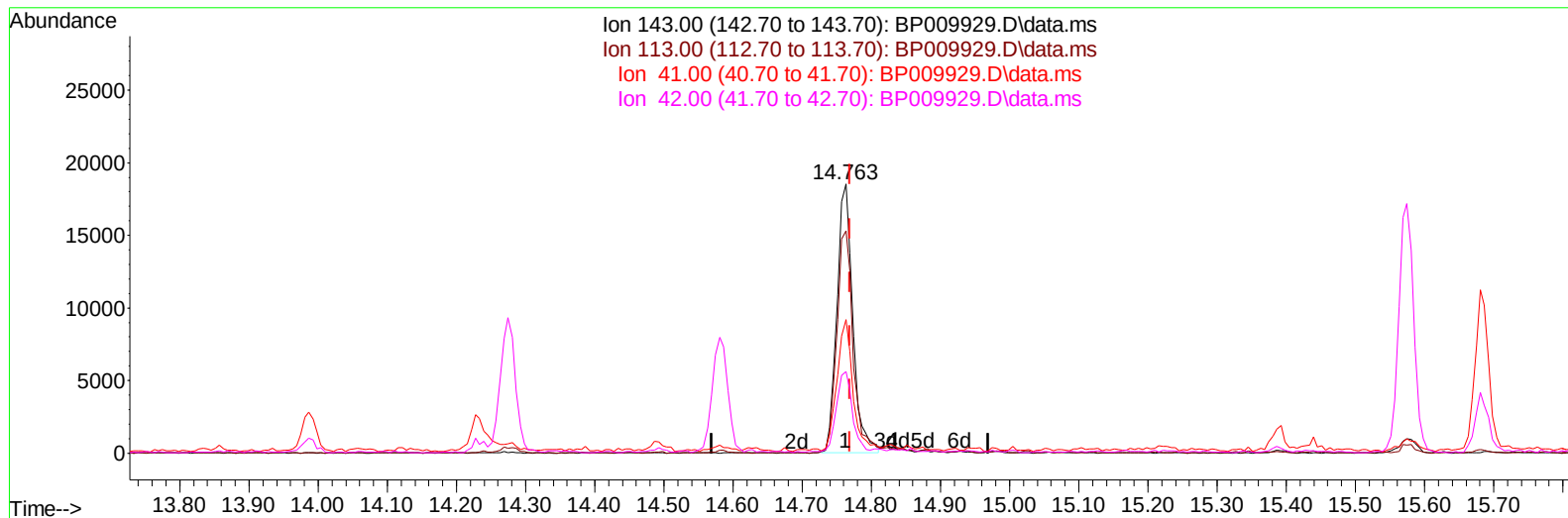
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TIC: BP009929.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.50 ng/u1

response 29503

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.69#
41.00	23.20	49.63#
42.00	18.00	30.31#

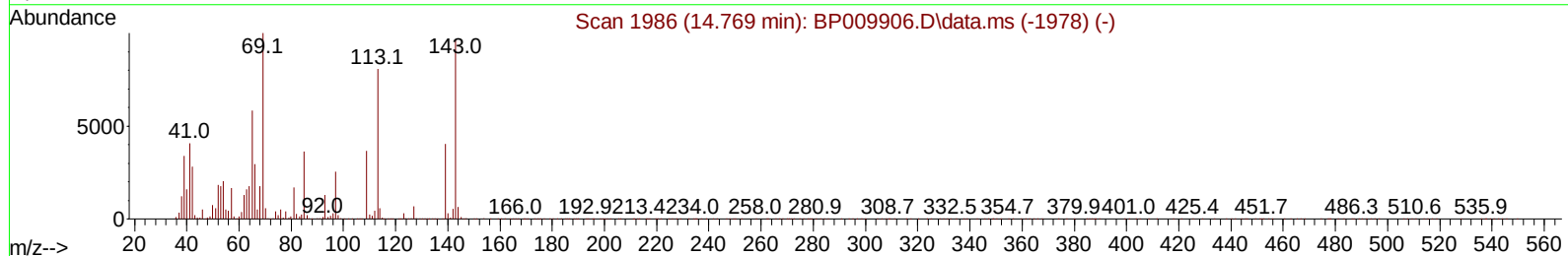
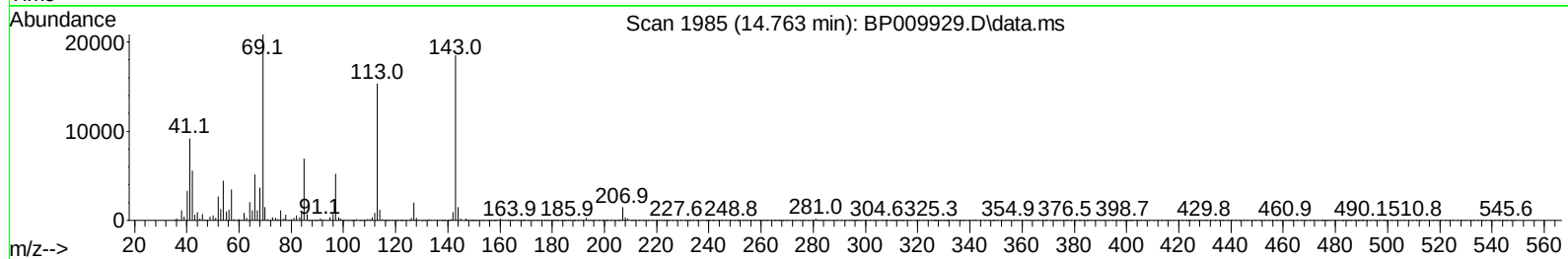
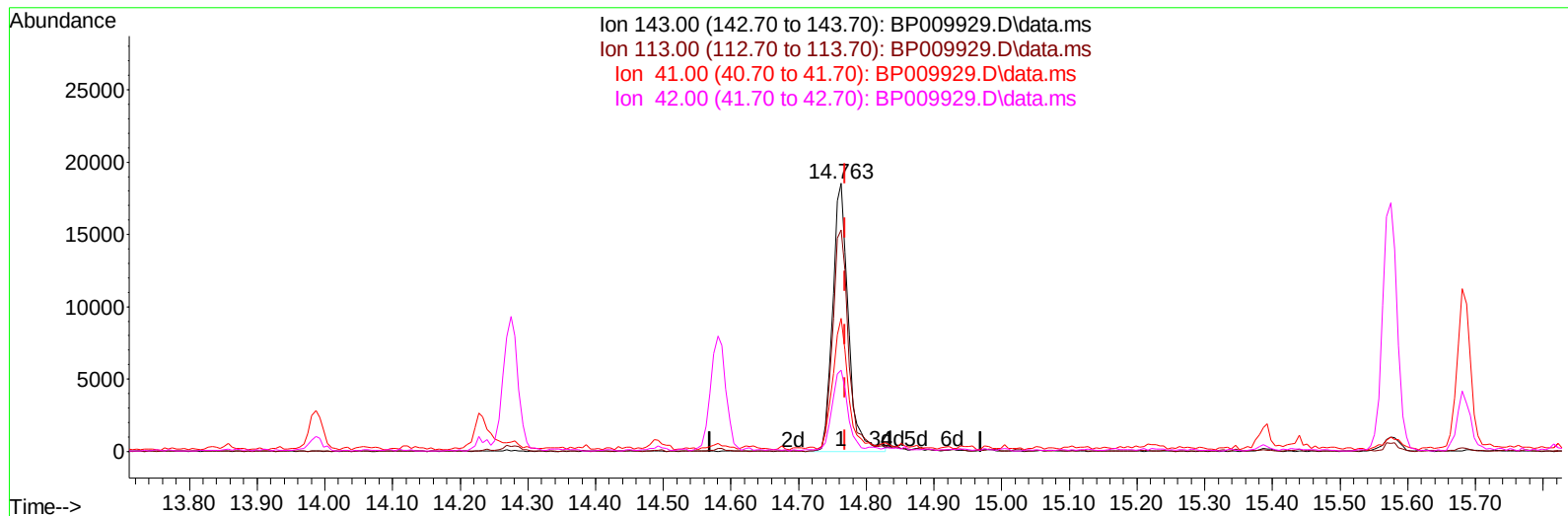
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Sample : N2421-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J56

Manual IntegrationsAPPROVED

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TIC: BP009929.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.61 ng/ul m

response 30071

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.69#
41.00	23.20	49.63#
42.00	18.00	30.31#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	119403	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	526535	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.581	164	376349	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.333	188	834117	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	846781	20.000	ng/ul	0.00
88) Perylene-d12	23.874	264	759570	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12996m	4.858	ng/uL	0.00
4) Pyridine-d5	3.799	84	60860	8.251	ng/ul	0.00
7) Phenol-d5	7.105	99	70409	6.717	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	207458	33.242	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	197059	24.921	ng/ul	0.00
15) 4-Methylphenol-d8	8.651	113	134708	15.394	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	134635	35.458	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	139149	33.426	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	240014	27.553	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	300974	24.259	ng/ul	0.00
46) Dimethylphthalate-d6	13.986	166	1000943	33.858	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1100849	31.432	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	30071m	5.609	ng/ul	0.00
60) Fluorene-d10	15.575	176	909007	36.551	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	167261	33.849	ng/ul	0.00
73) Anthracene-d10	17.433	188	1622785	40.634	ng/ul	0.00
81) Pyrene-d10	19.663	212	2004197	43.337	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	1668023	42.169	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.134	94	13885	1.323	ng/ul#	83

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

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