

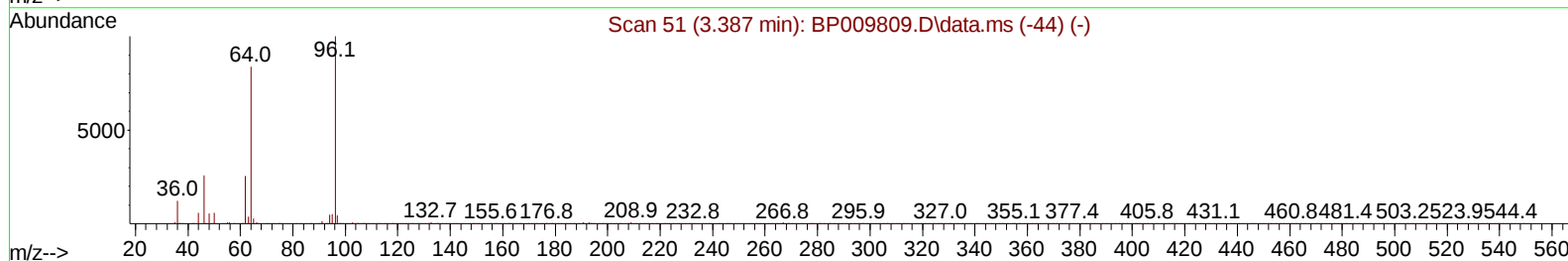
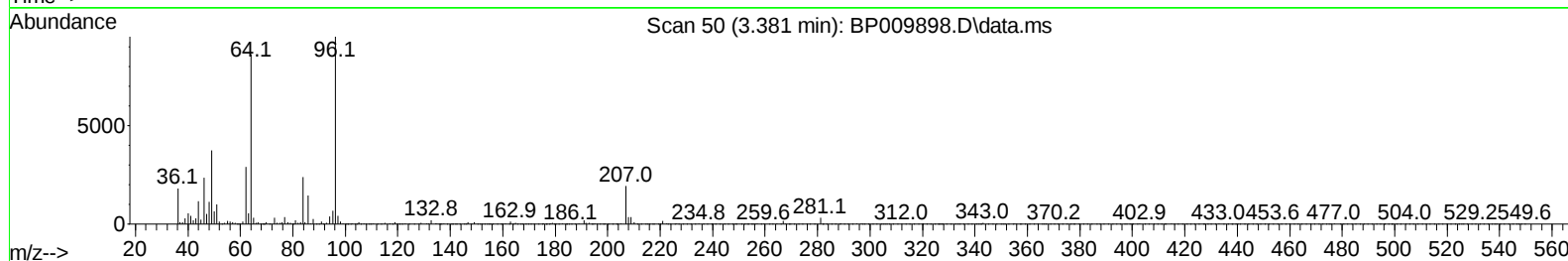
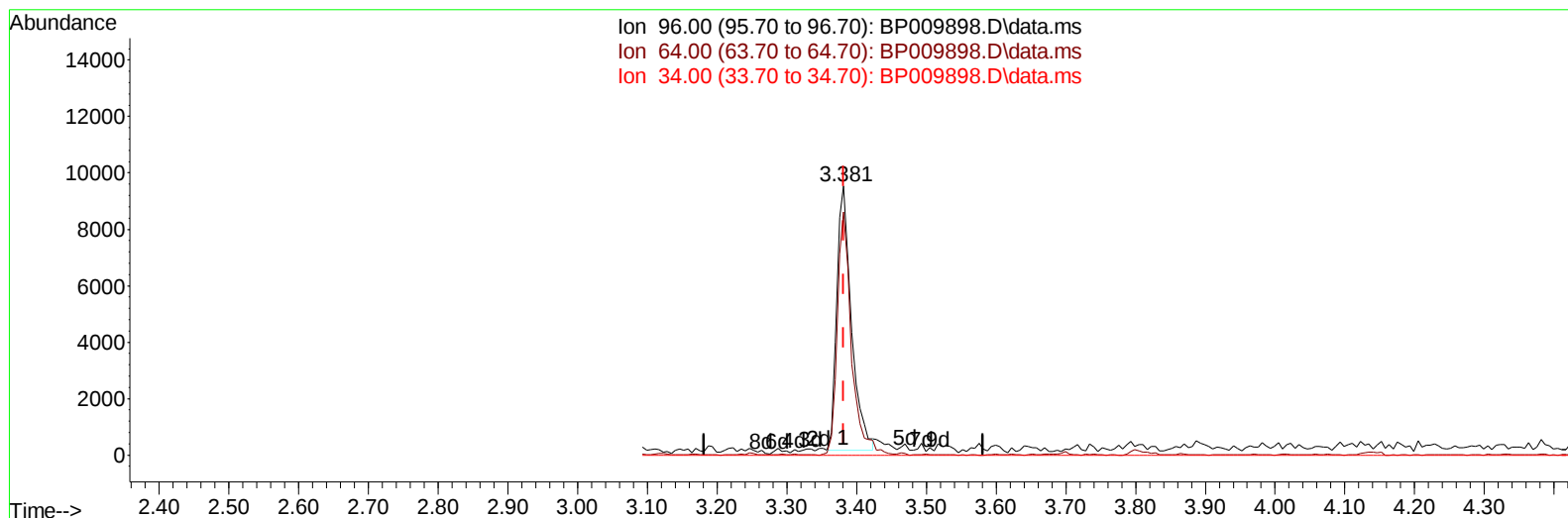
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009898.D
 Acq On : 16 Apr 2022 12:55
 Operator : CG/JU
 Sample : N2421-17
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J77

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:12:57 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 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022



TIC: BP009898.D\data.ms

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

3.381min (+ 0.000) 4.42 ng/uL

response 13554

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	90.48#
34.00	0.00	0.00
0.00	0.00	0.00

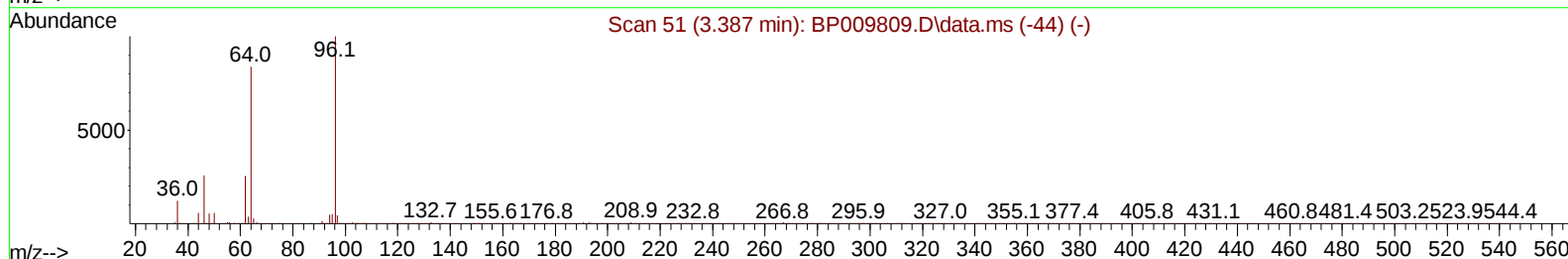
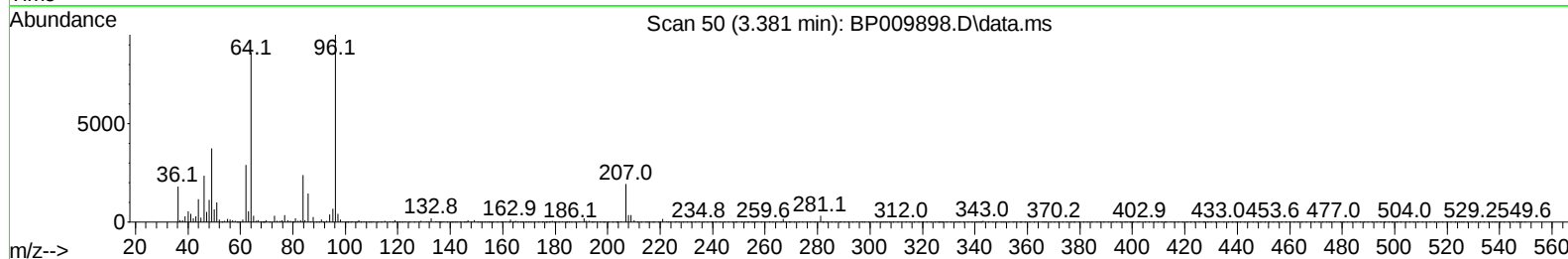
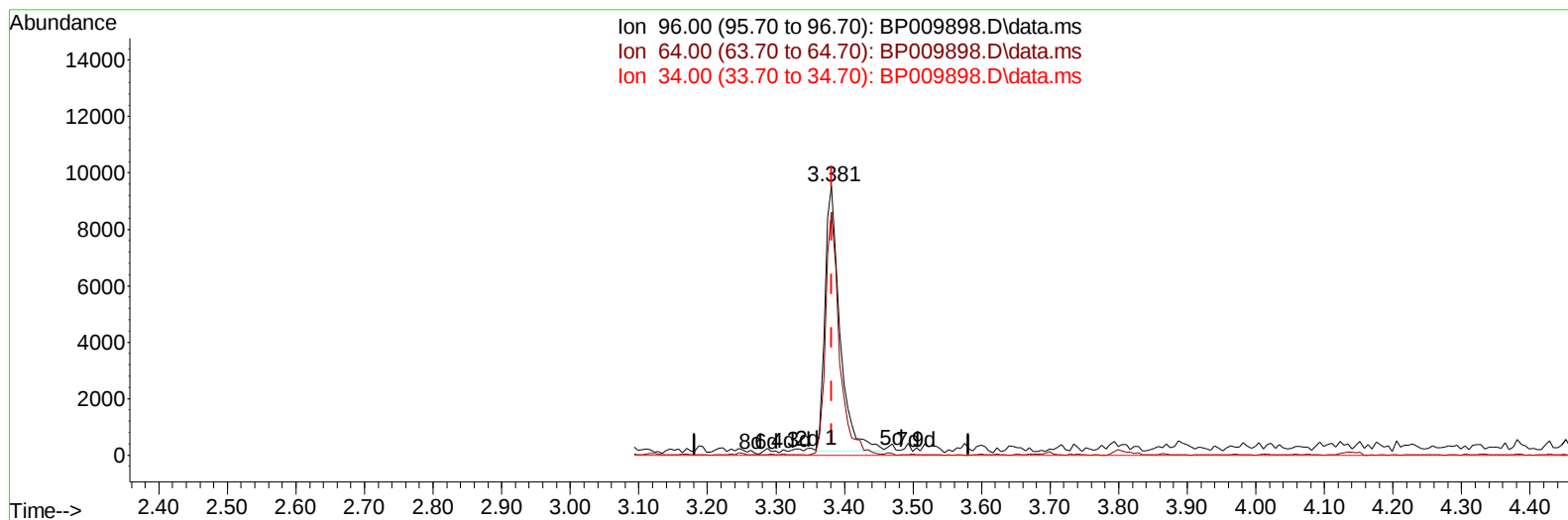
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TIC: BP009898.D\data.ms

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

3.381min (+ 0.000) 4.62 ng/uL m

response 14178

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	90.48#
34.00	0.00	0.00
0.00	0.00	0.00

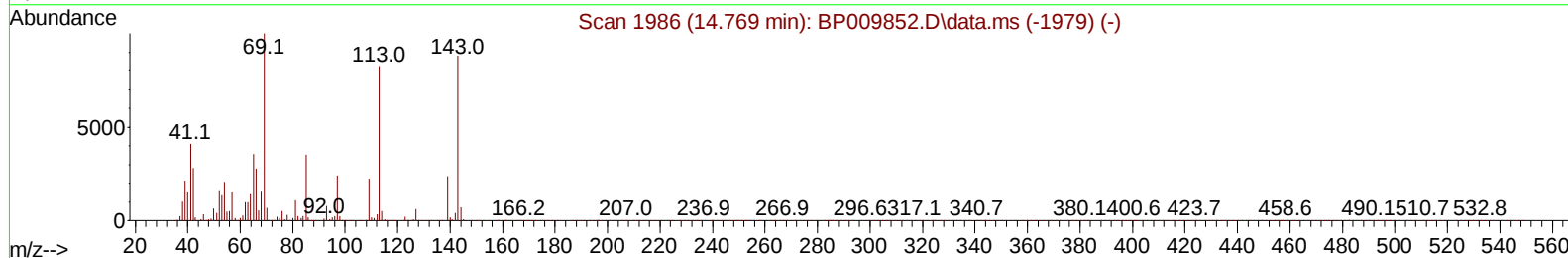
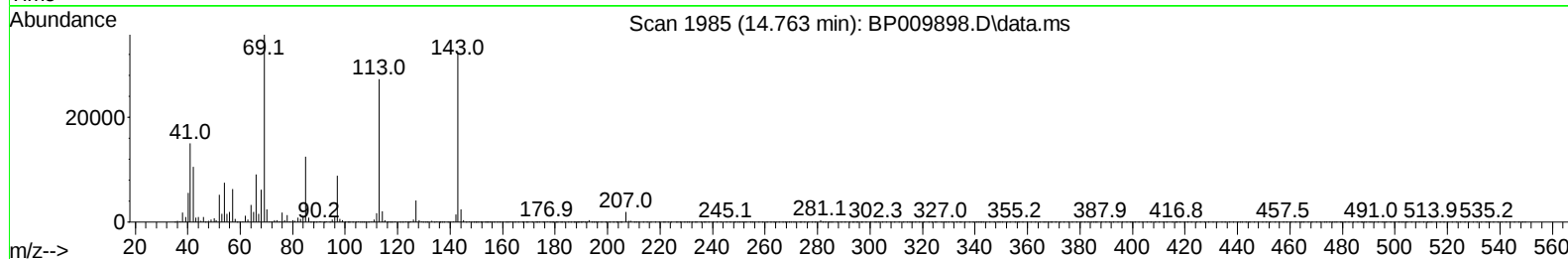
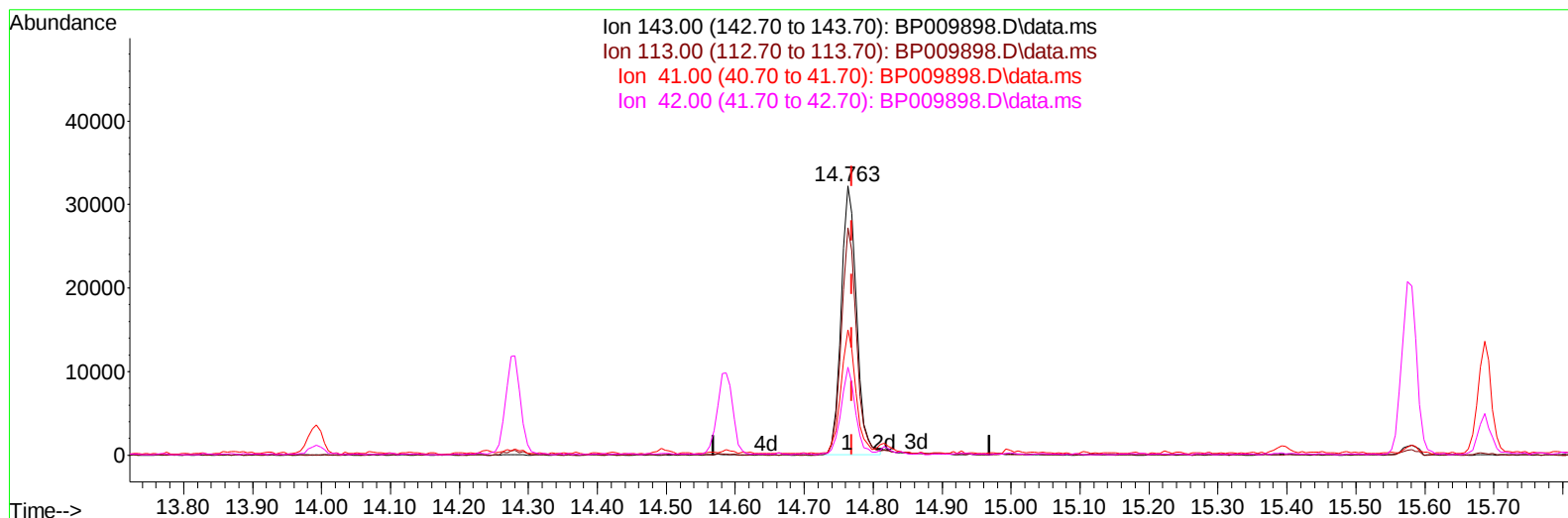
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TIC: BP009898.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 7.29 ng/u1

response 49790

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.52#
41.00	23.20	46.59#
42.00	18.00	32.71#

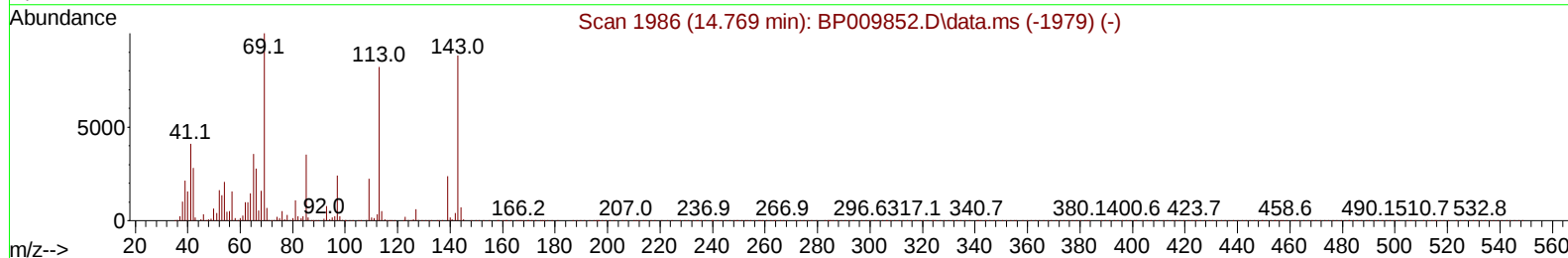
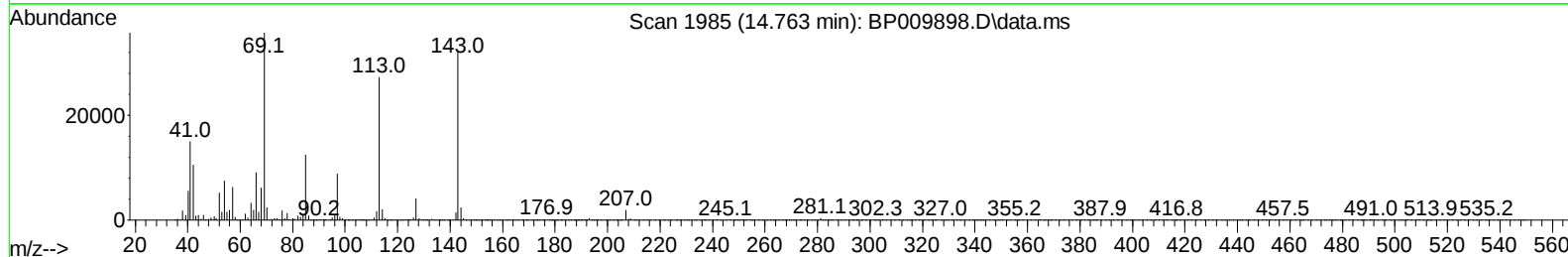
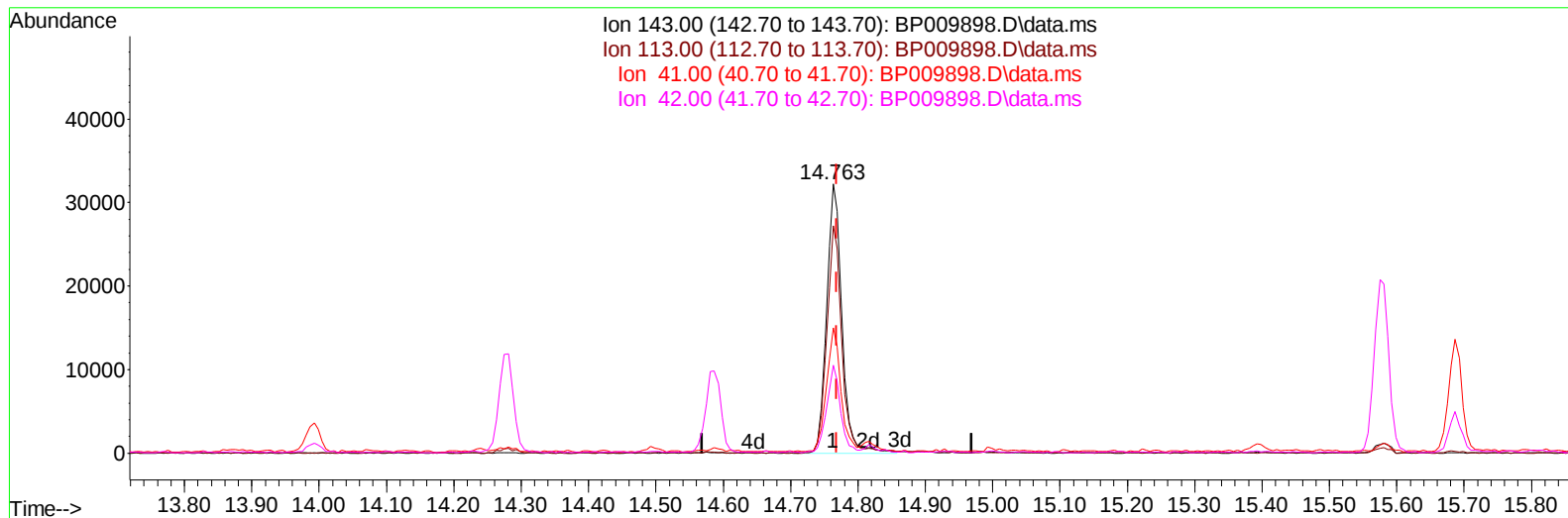
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Operator : CG/JU
Sample : N2421-17
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP009898.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 7.48 ng/ul m

response 51071

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.52#
41.00	23.20	46.59#
42.00	18.00	32.71#

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 ALS Vial : 50 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	136920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	638874	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	479299	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1079009	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	1046486	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	927886	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	14178m	4.621	ng/uL	0.00
4) Pyridine-d5	3.799	84	82726	9.780	ng/ul	0.00
7) Phenol-d5	7.111	99	96412	8.020	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	246170	34.399	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	245041	27.024	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	192161	19.150	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	166874	36.221	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	180369	35.710	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	311324	29.455	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	421322	27.987	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1307018	34.715	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1458721	32.704	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	51071m	7.480	ng/ul	0.00
60) Fluorene-d10	15.581	176	1147709	36.237	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	210956	33.003	ng/ul	0.00
73) Anthracene-d10	17.439	188	2047690	39.636	ng/ul	0.00
81) Pyrene-d10	19.669	212	2481438	43.417	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	2075123	42.945	ng/ul	0.00

Target Compounds Qvalue

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

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