

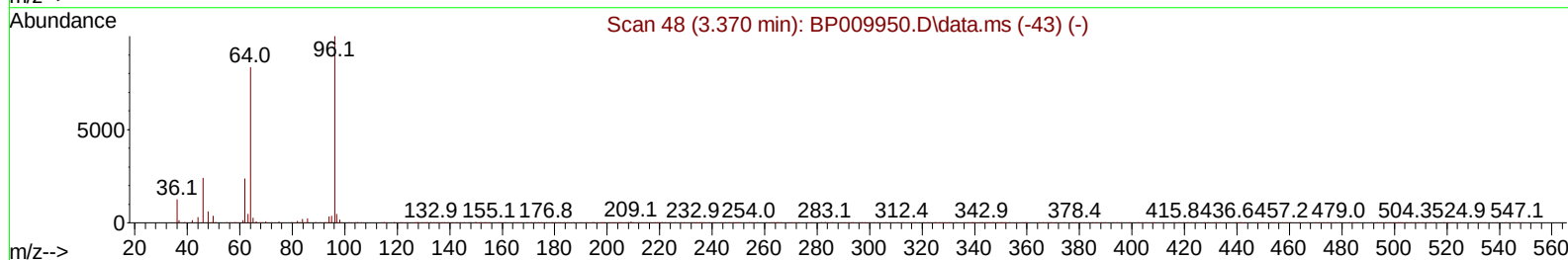
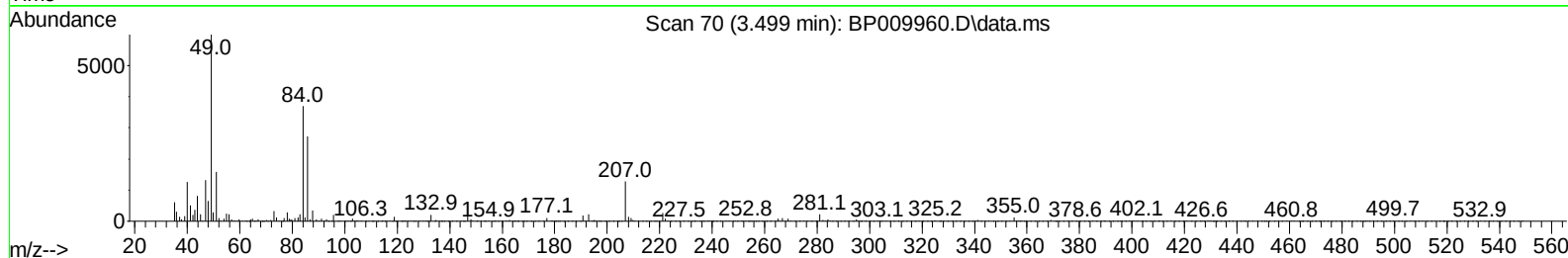
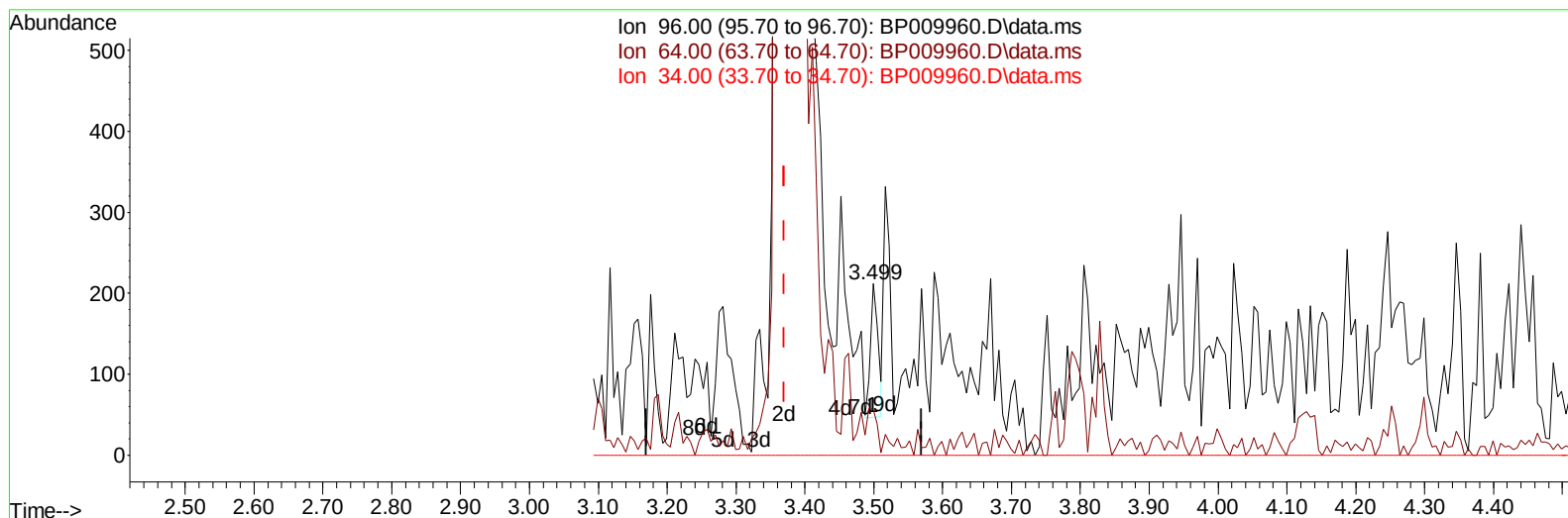
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009960.D
 Acq On : 19 Apr 2022 23:40
 Operator : CG/JU
 Sample : N2422-17
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J12

Manual IntegrationsAPPROVED

Quant Time: Apr 20 01:34:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/20/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP009960.D\data.ms

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

3.499min (+ 0.129) 0.05 ng/uL

response 130

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	63.68
34.00	0.00	0.00
0.00	0.00	0.00

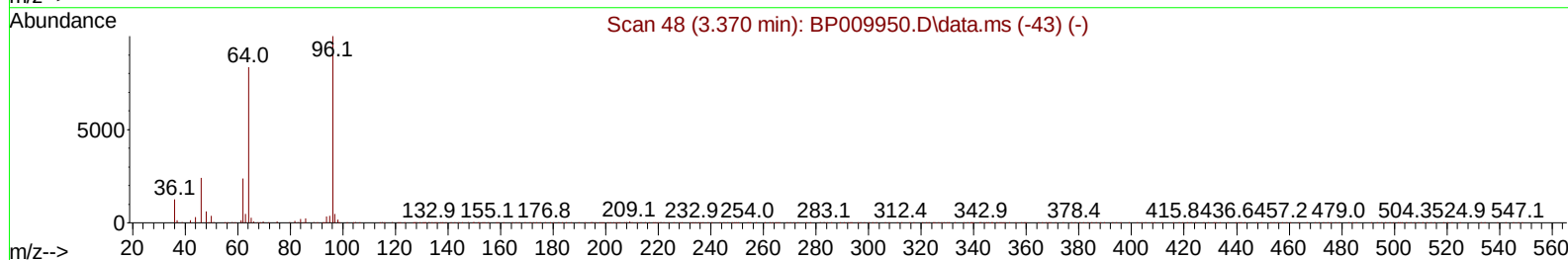
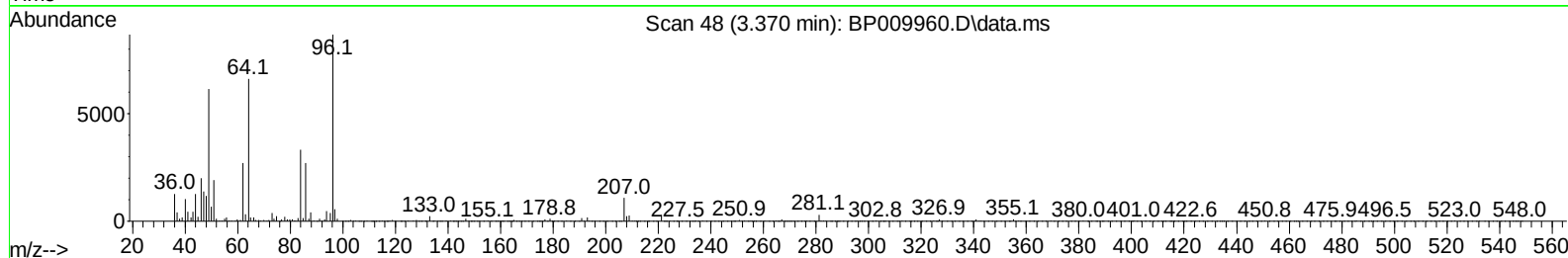
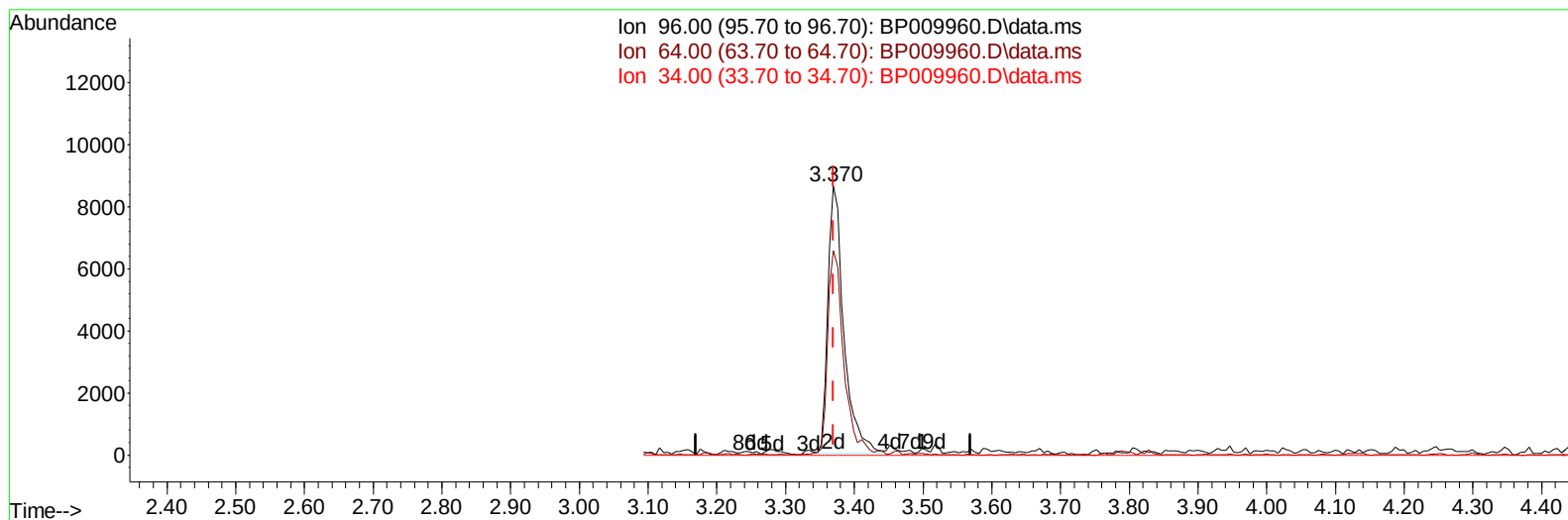
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TIC: BP009960.D\data.ms

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

3.370min (-0.000) 4.99 ng/uL m

response 13859

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	76.14#
34.00	0.00	0.00
0.00	0.00	0.00

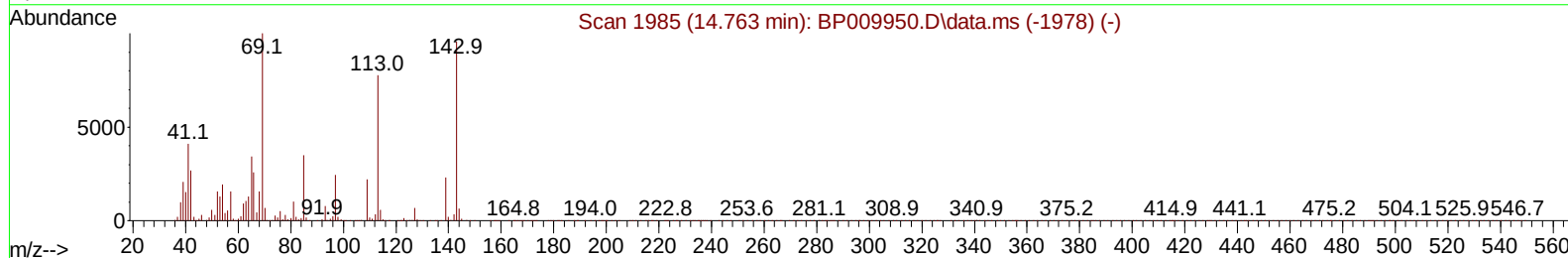
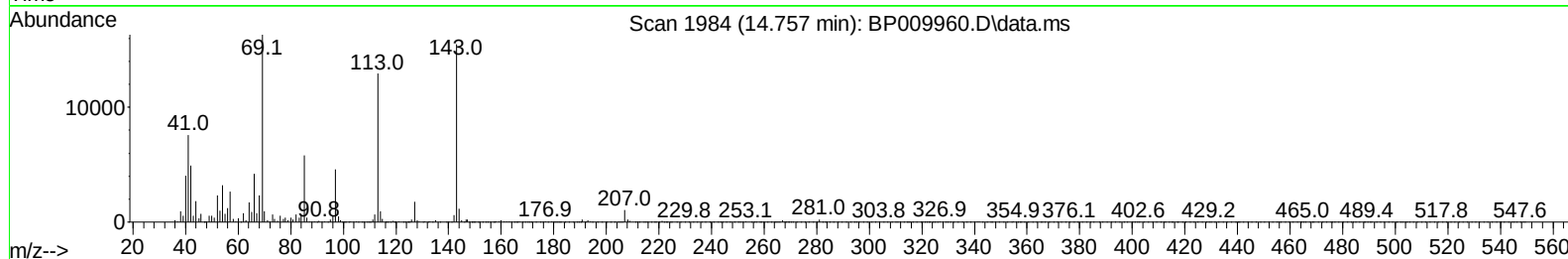
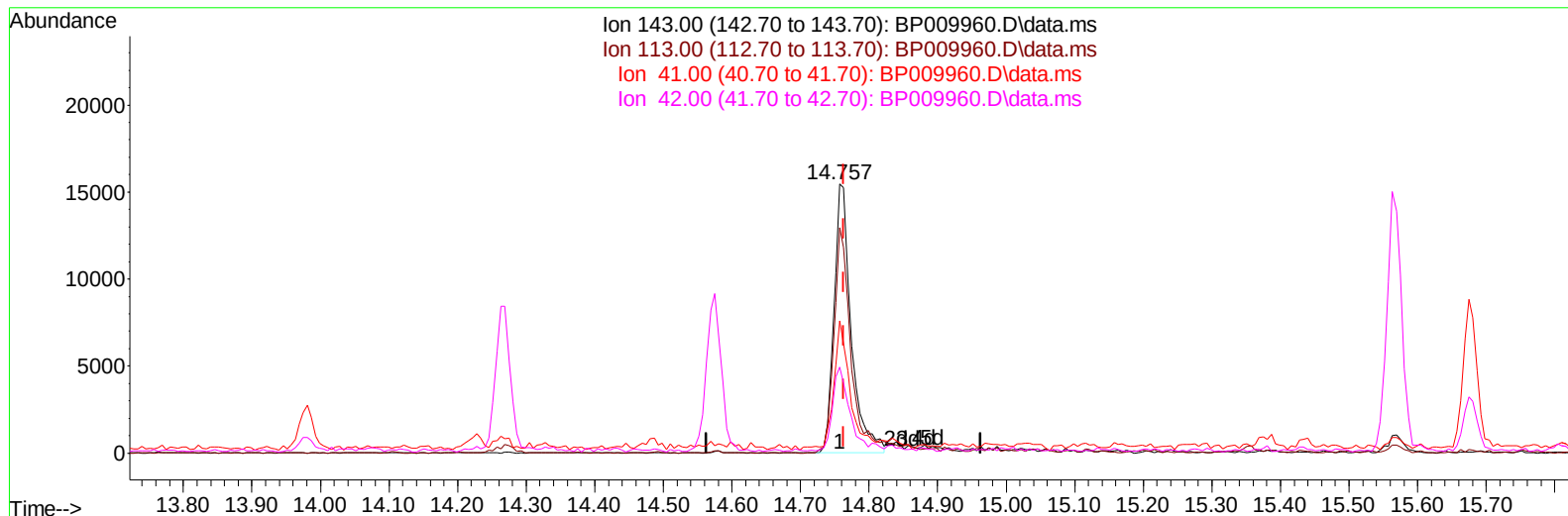
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP009960.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 4.67 ng/u1

response 26914

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.71#
41.00	23.20	49.18#
42.00	18.00	31.93#

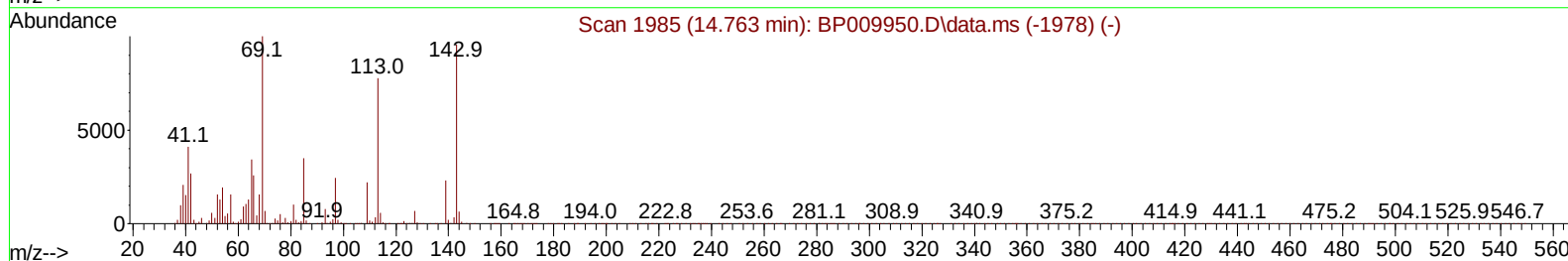
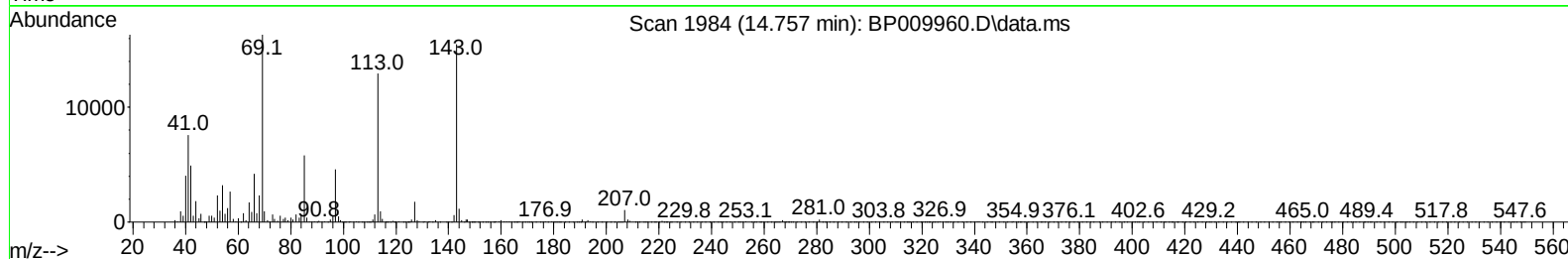
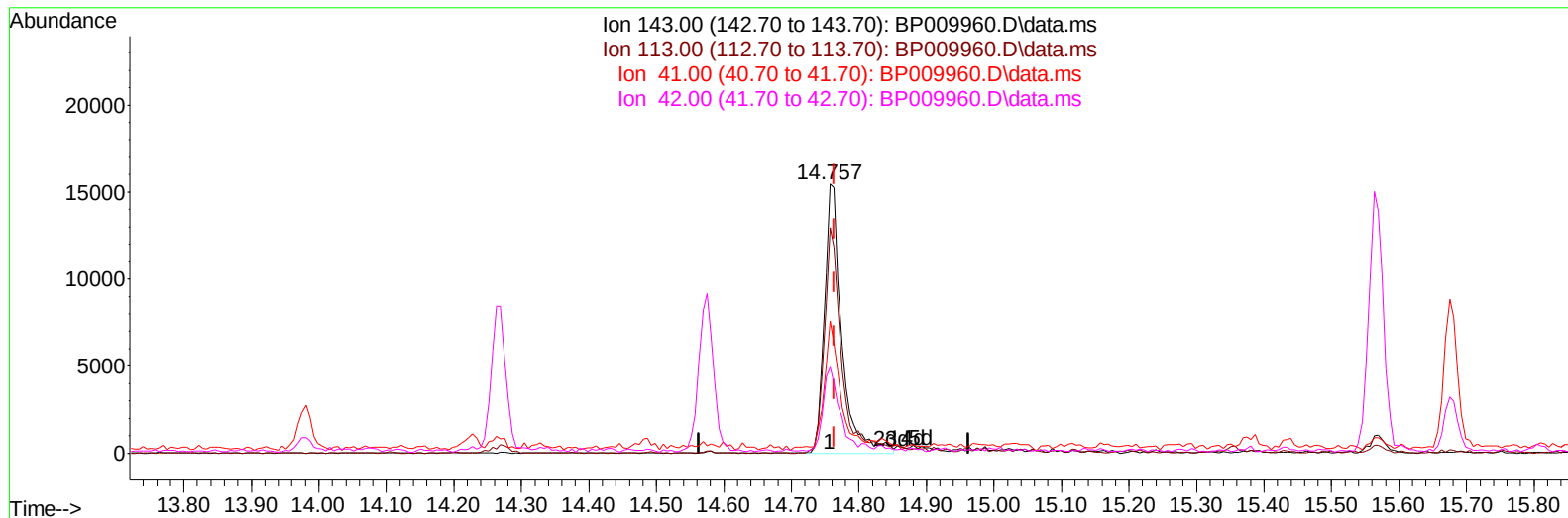
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Data File : BP009960.D
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Operator : CG/JU
Sample : N2422-17
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP009960.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 4.83 ng/ul m

response 27853

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.71#
41.00	23.20	49.18#
42.00	18.00	31.93#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	123999	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	551402	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	404796	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	906068	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	936284	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	922828	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	13859m	4.988	ng/uL	0.00
4) Pyridine-d5	3.793	84	57462	7.501	ng/ul	0.00
7) Phenol-d5	7.099	99	63340	5.818	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	181229	27.963	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	172301	20.982	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	126641	13.935	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	113510	28.546	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	120492	27.639	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	209180	22.930	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	297714	22.914	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	939238	29.538	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	996905	26.464	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	27853m	4.830	ng/ul	0.00
60) Fluorene-d10	15.569	176	819795	30.648	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	138361	25.777	ng/ul	0.00
73) Anthracene-d10	17.428	188	1403398	32.350	ng/ul	0.00
81) Pyrene-d10	19.657	212	1718915	33.615	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1561640	32.495	ng/ul	0.00

Target Compounds Qvalue

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

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