

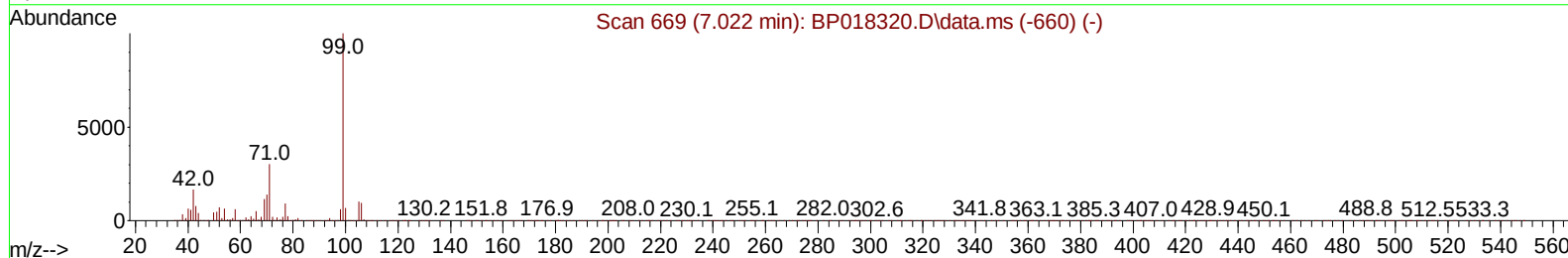
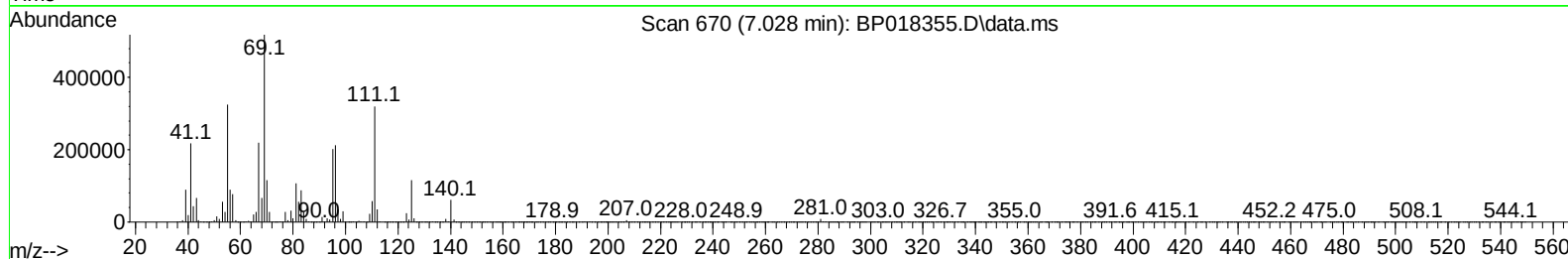
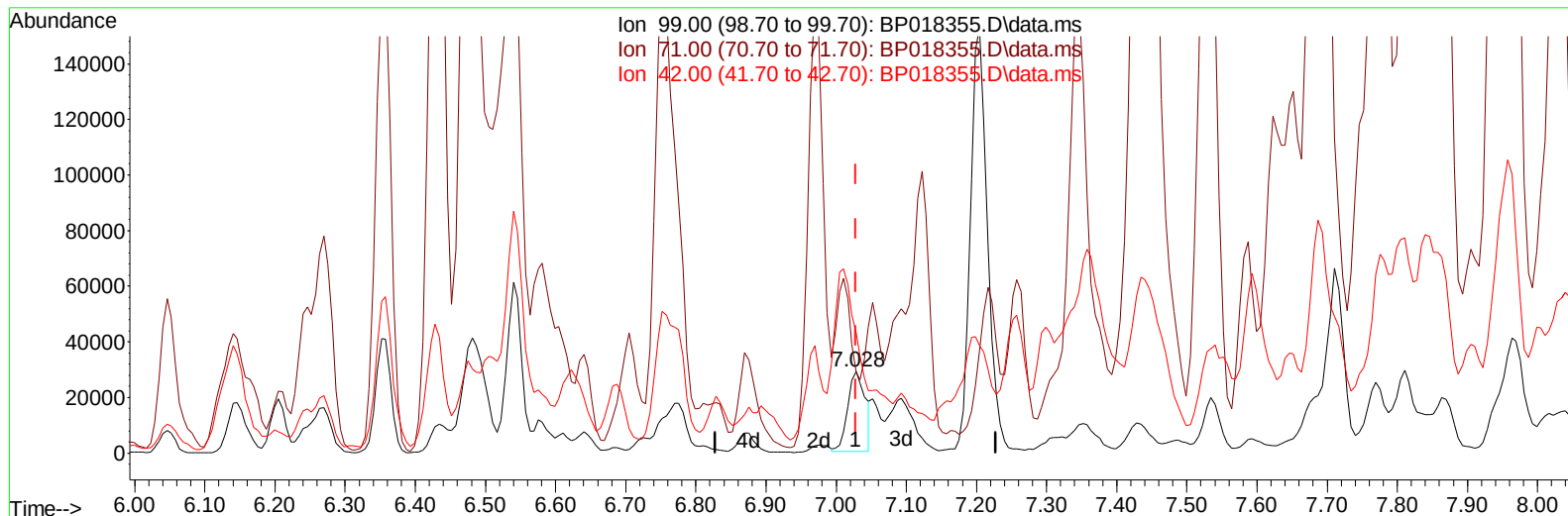
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018355.D
 Acq On : 25 Nov 2023 07:27
 Operator : MA/JU
 Sample : 05264-18
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKS9

Manual IntegrationsAPPROVED

Quant Time: Nov 27 00:17:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/27/2023
 Supervised By :mohammad ahmed 11/27/2023



TIC: BP018355.D\data.ms

(7) Phenol-d5 (S)

7.028min (+ 0.000) 1.79 ng/ul

response 50493

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	31.30	98.96#
42.00	17.40	147.37#
0.00	0.00	0.00

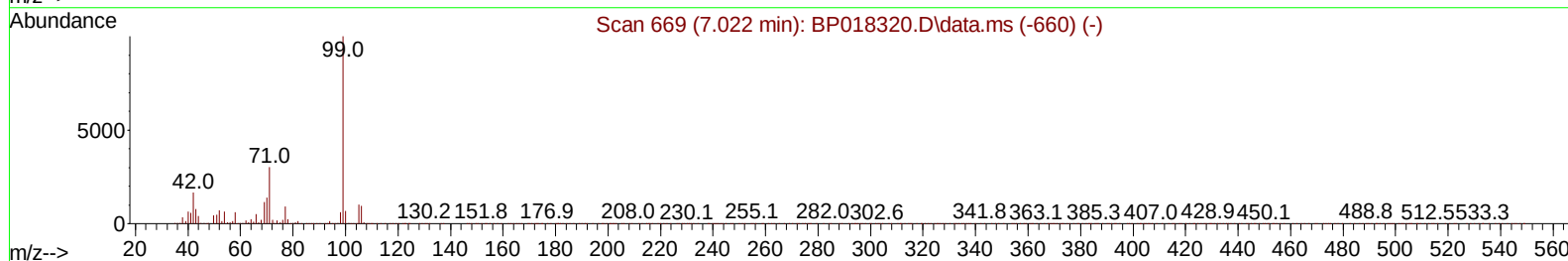
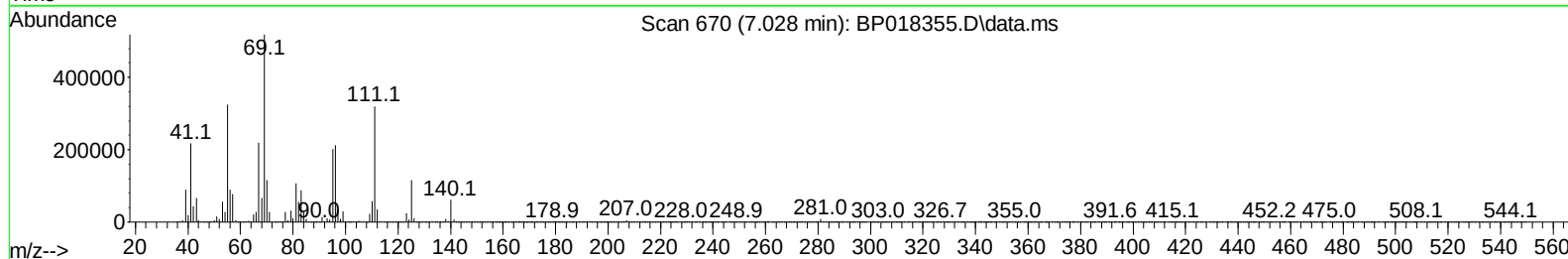
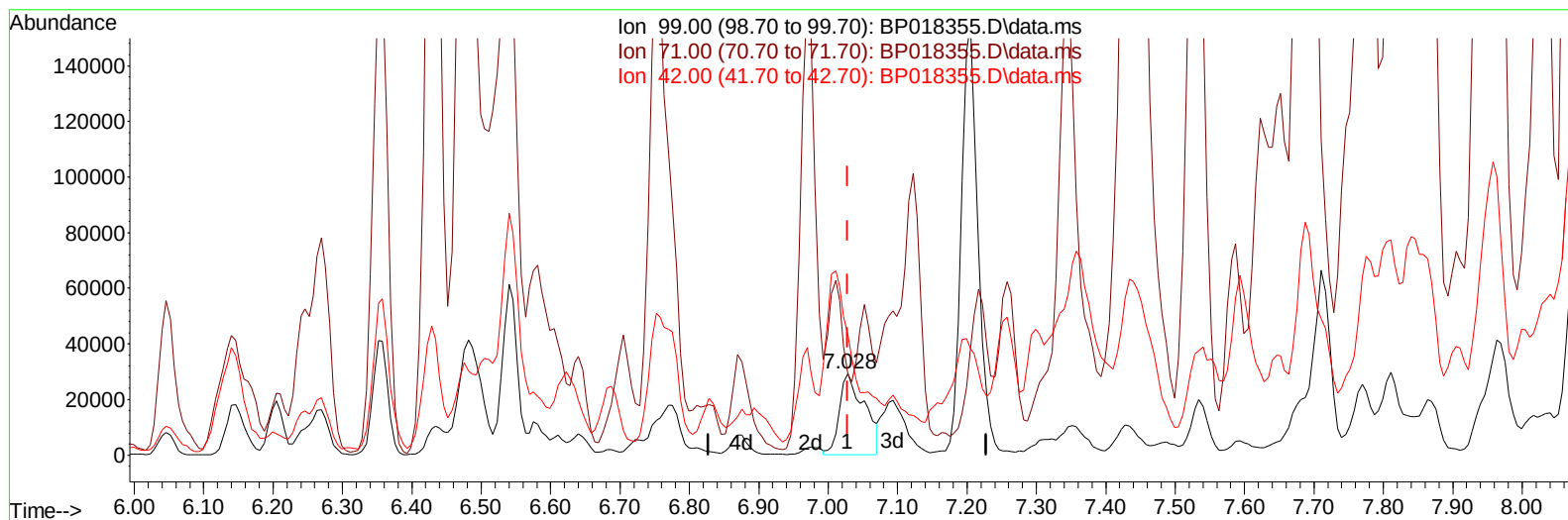
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TIC: BP018355.D\data.ms

(7) Phenol-d5 (S)

7.028min (+ 0.000) 2.61 ng/ul m

response 73427

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	31.30	98.96#
42.00	17.40	147.37#
0.00	0.00	0.00

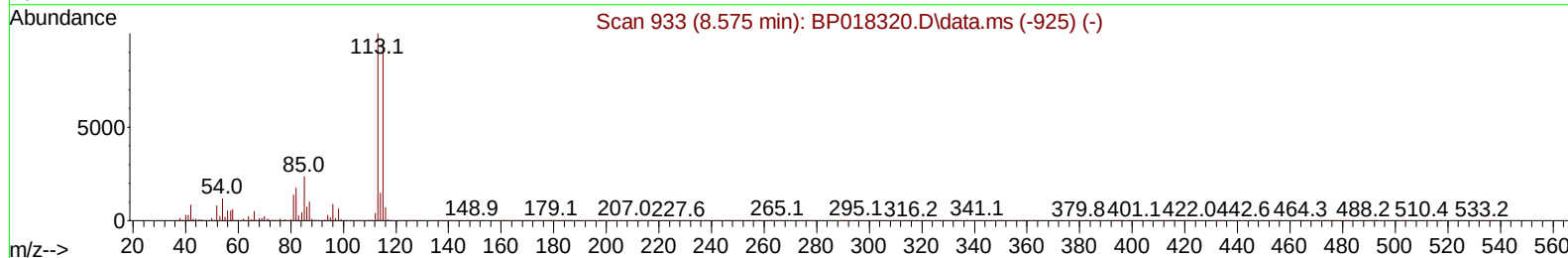
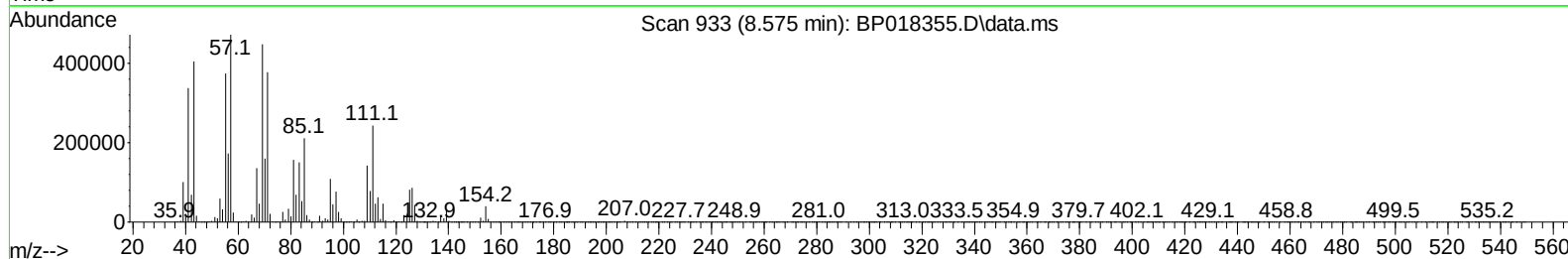
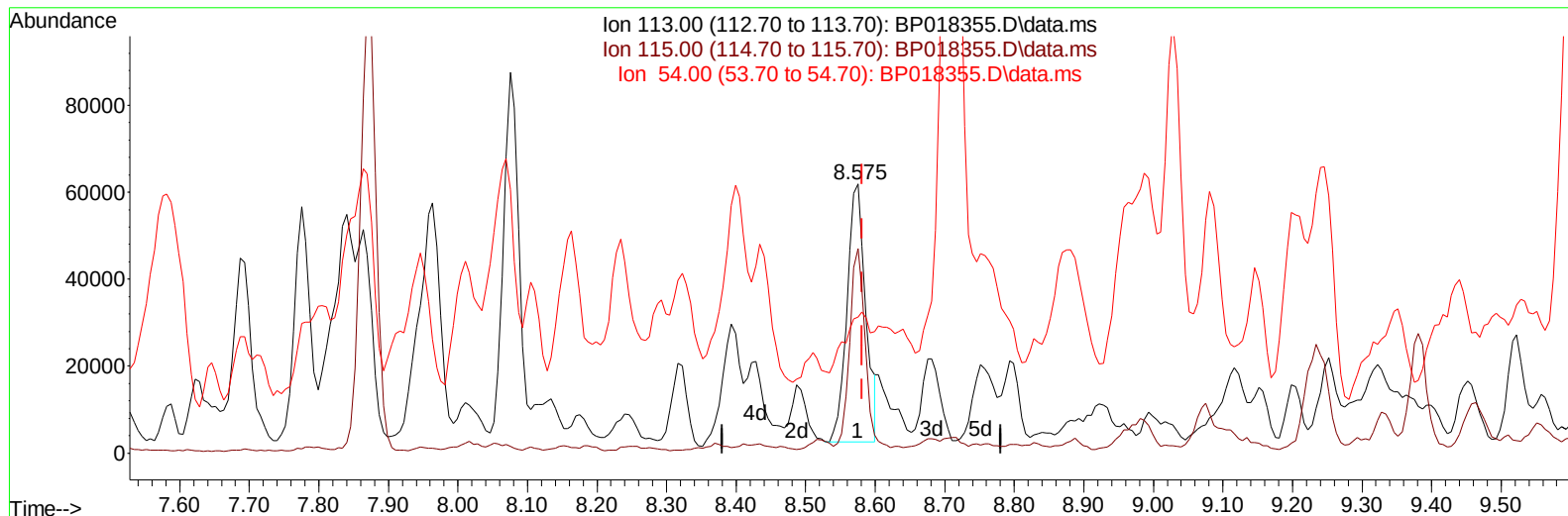
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Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Nov 27 00:18:12 2023
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TIC: BP018355.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.575min (-0.006) 4.88 ng/uI

response 110906

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	93.40	75.93
54.00	12.90	50.55#
0.00	0.00	0.00

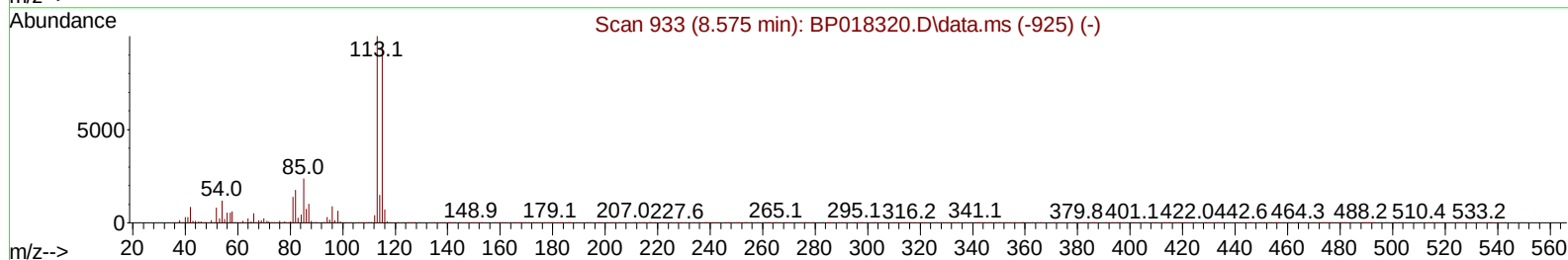
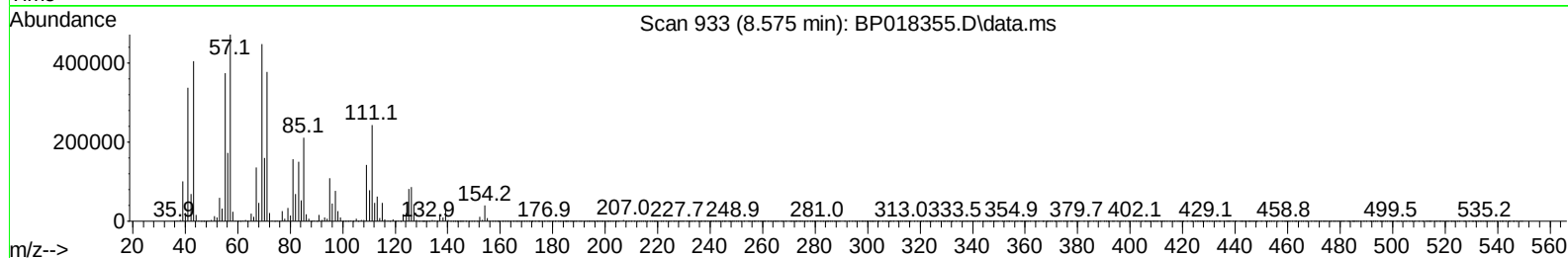
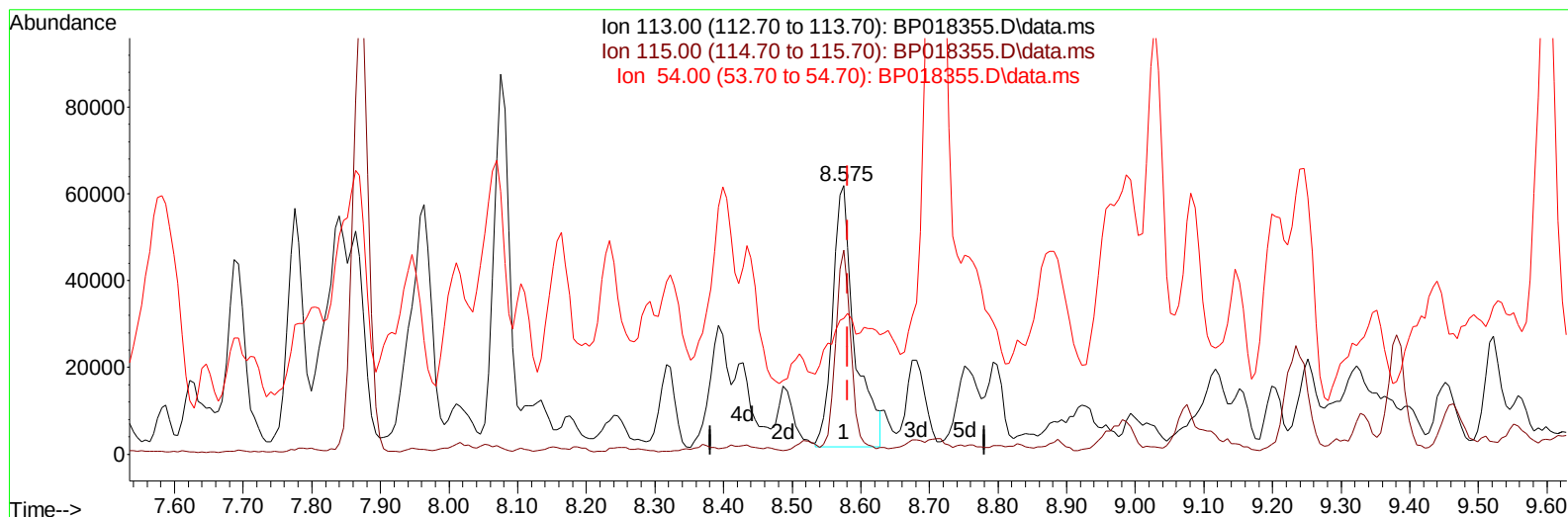
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Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKS9

Manual IntegrationsAPPROVED

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TIC: BP018355.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.575min (-0.006) 5.92 ng/ul m

response 134530

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	93.40	75.93
54.00	12.90	50.55#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	307905	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	1142270	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.498	164	722240	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1535465	20.000	ng/ul	0.00
79) Chrysene-d12	21.357	240	1464911	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1794877	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	23702	2.883	ng/uL	0.00
4) Pyridine-d5	3.693	84	77278	3.421	ng/ul	0.00
7) Phenol-d5	7.028	99	73427m	2.609	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	316108	18.786	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	98833	4.183	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	134530m	5.924	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	96936	10.448	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	39979	4.666	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	108347	5.241	ng/ul	0.00
31) 4-Chloroaniline-d4	10.805	131	101076	3.625	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	731972	11.390	ng/ul	0.00
49) Acenaphthylene-d8	14.193	160	808934	11.047	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	9193	1.036	ng/ul	0.00
60) Fluorene-d10	15.493	176	649317	12.080	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	31091	4.757	ng/ul	0.00
73) Anthracene-d10	17.345	188	1000452	12.190	ng/ul	0.00
81) Pyrene-d10	19.586	212	1284002	11.013	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	1235719	11.875	ng/ul	0.01
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.298	149	4354331	55.917	ng/ul#	Qvalue 96

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

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