

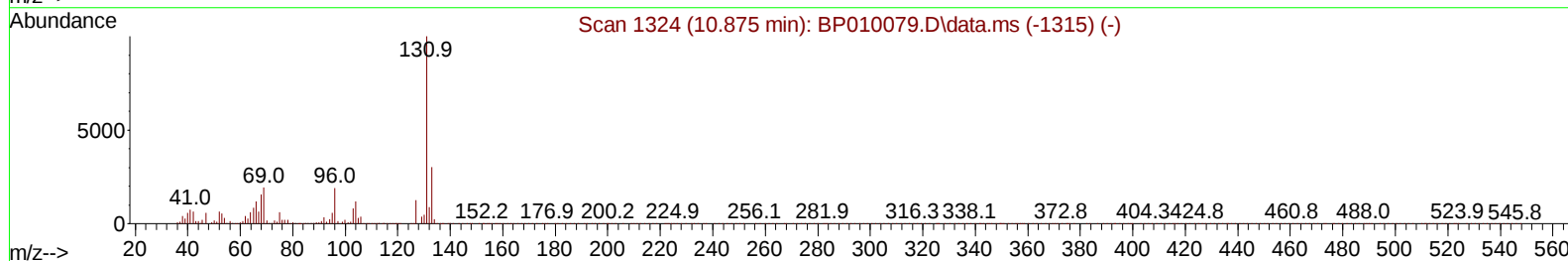
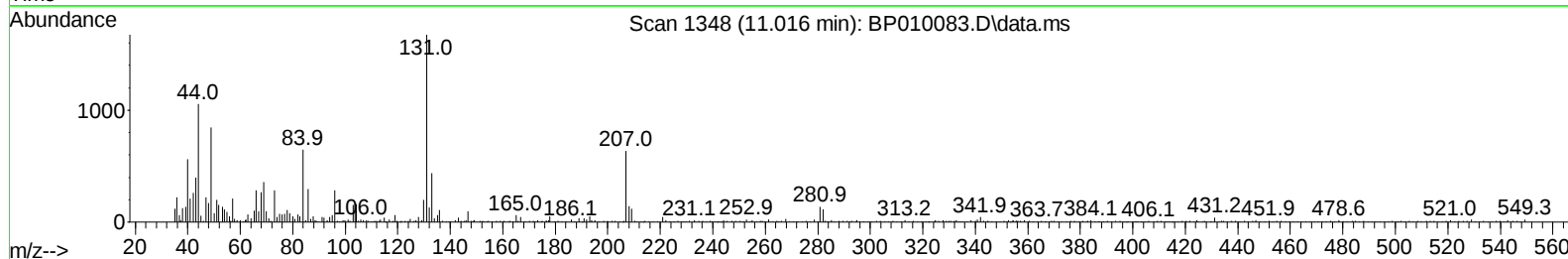
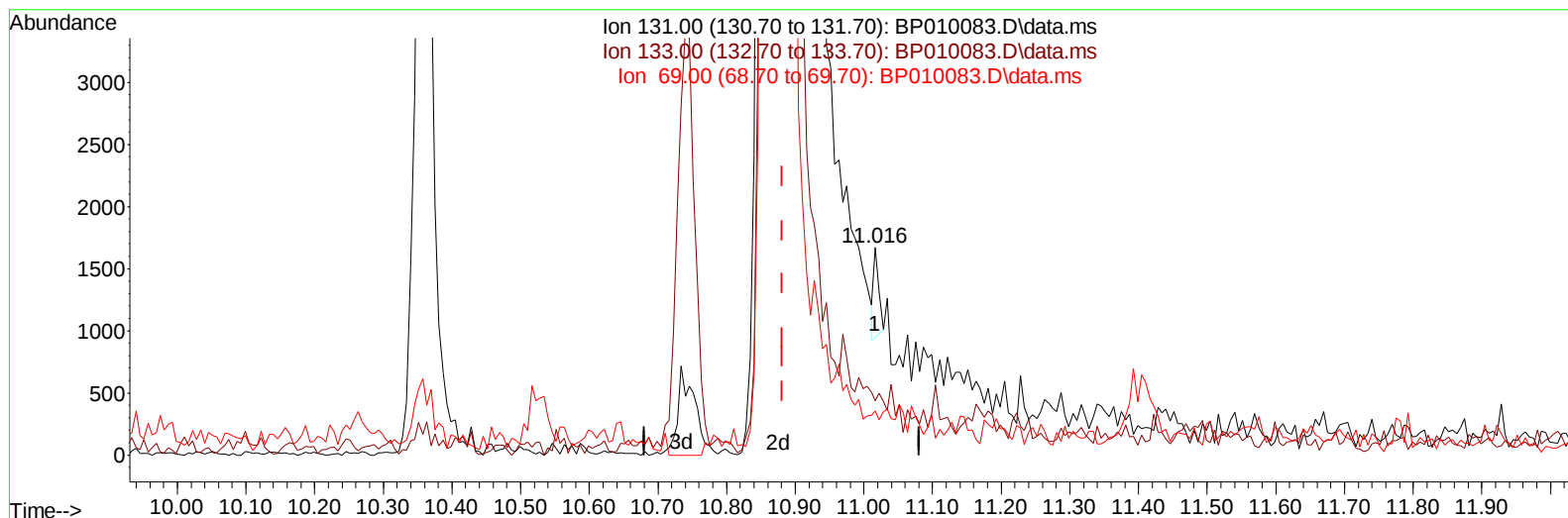
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010083.D
 Acq On : 25 Apr 2022 15:25
 Operator : CG/JU
 Sample : N2471-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFG2

Manual IntegrationsAPPROVED

Quant Time: Apr 26 01:02:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010083.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.016min (+ 0.135) 0.02 ng/u1

response 390

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	26.22
69.00	26.10	21.33
0.00	0.00	0.00

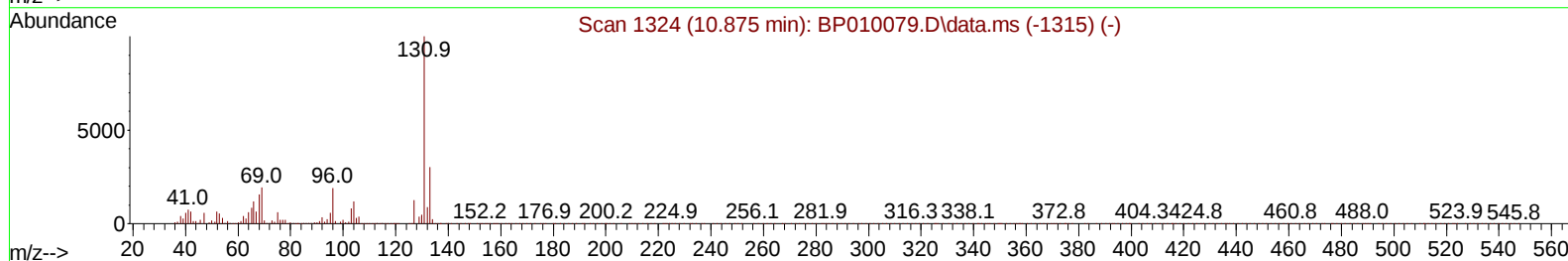
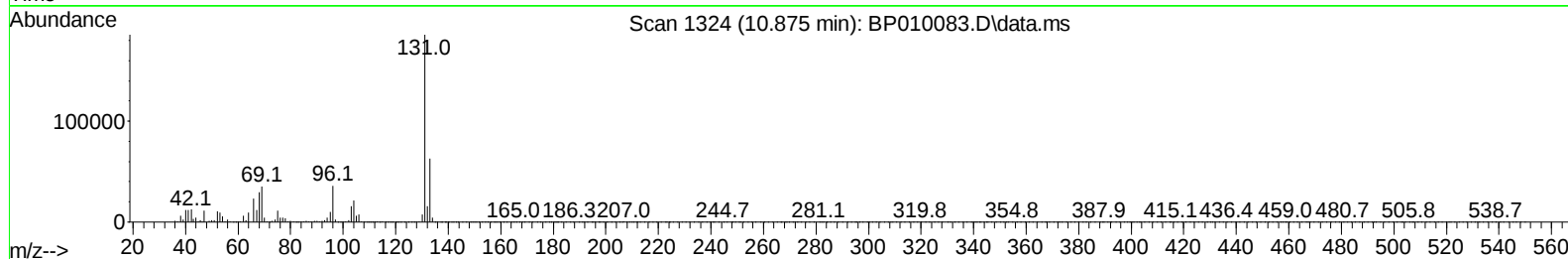
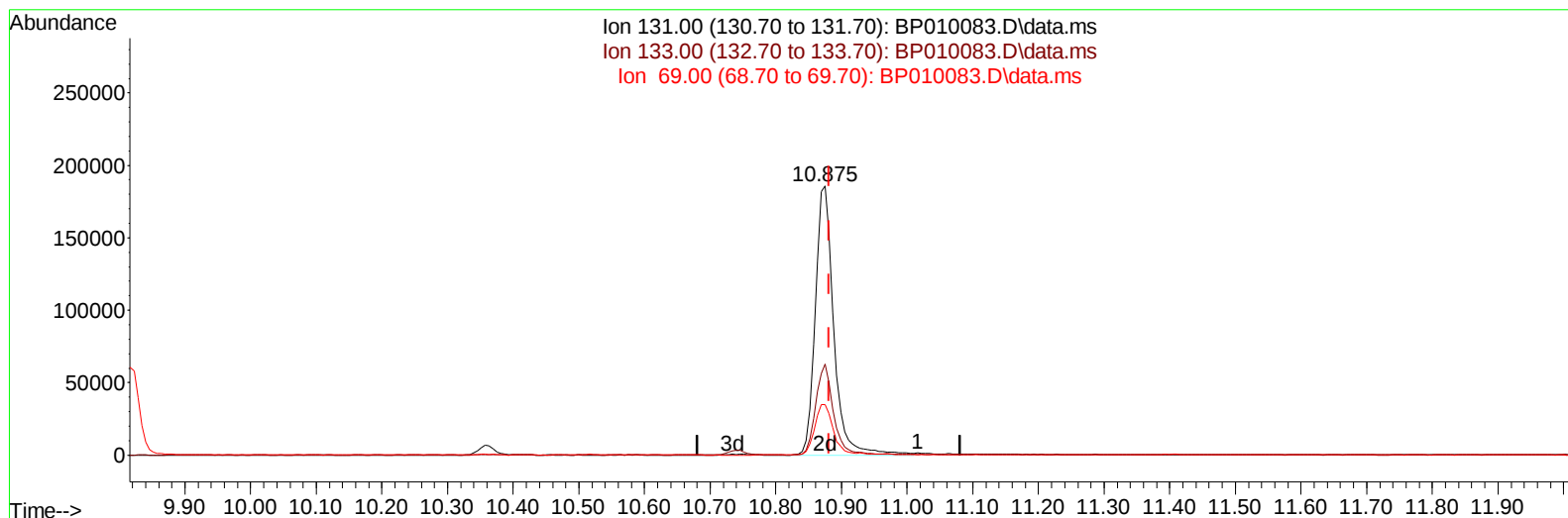
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TIC: BP010083.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.875min (-0.006) 20.86 ng/ul m

response 365893

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	33.91
69.00	26.10	18.87#
0.00	0.00	0.00

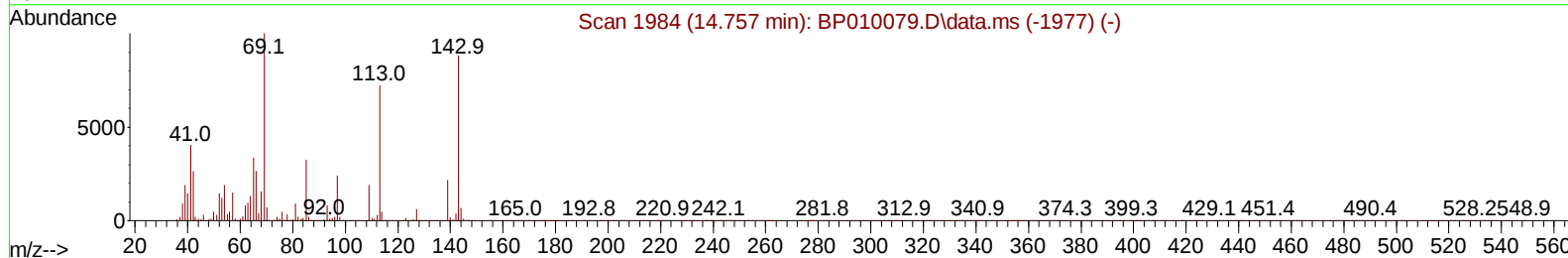
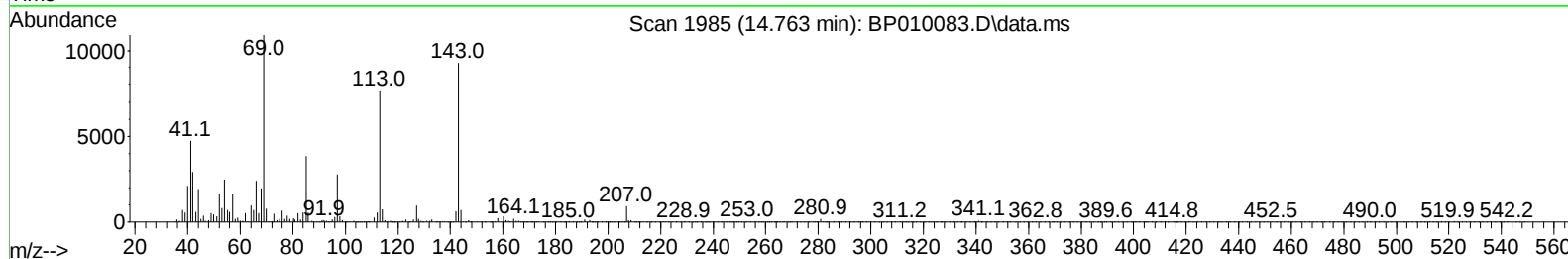
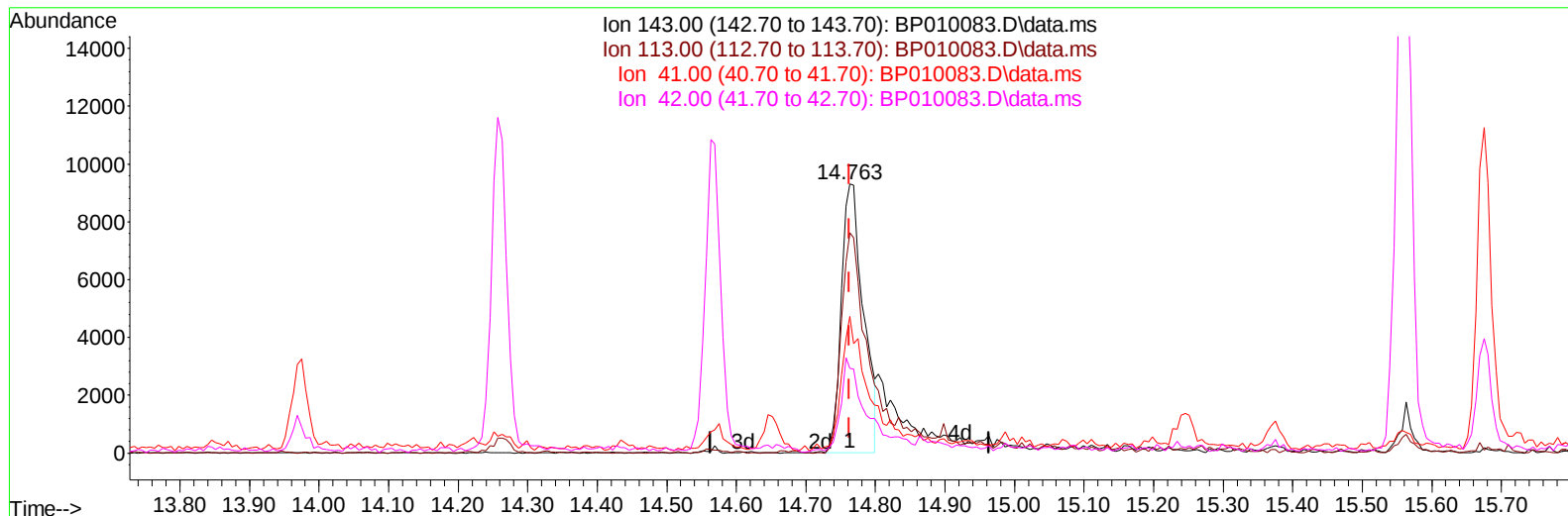
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 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP010083.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 2.85 ng/u1

response 20513

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.88#
41.00	23.20	50.82#
42.00	18.00	31.52#

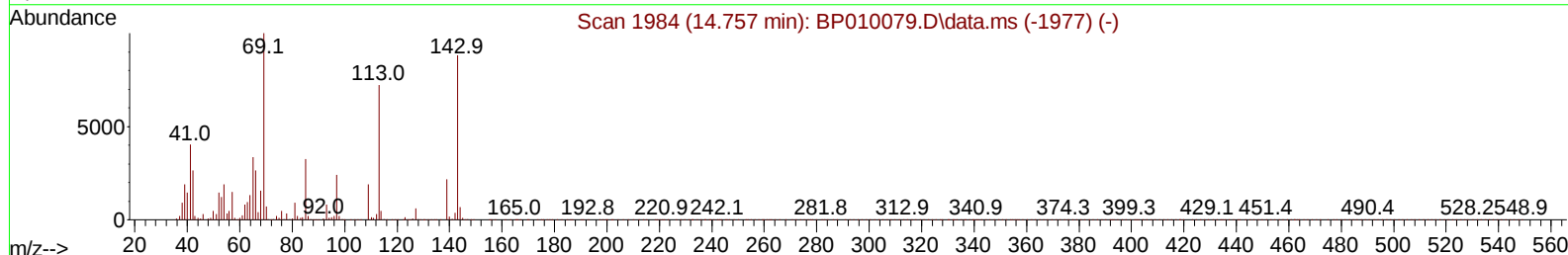
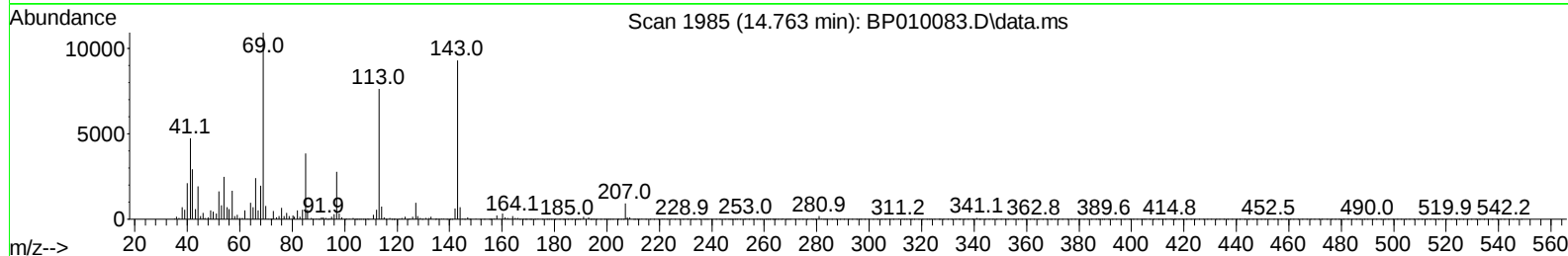
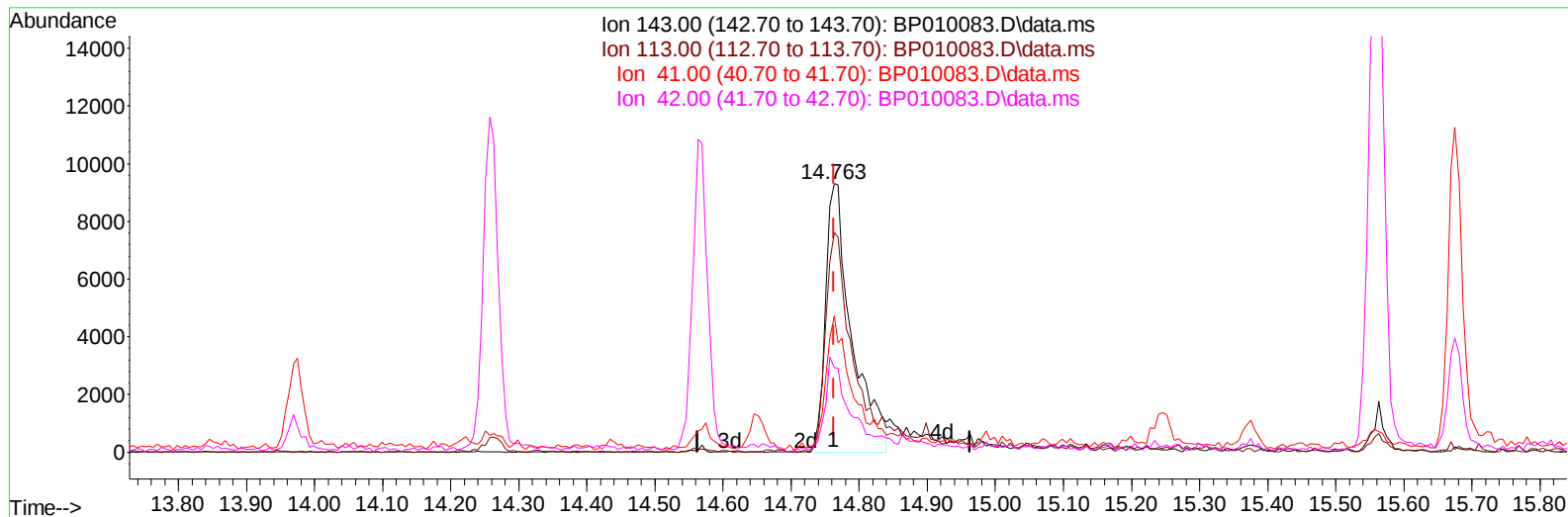
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 Operator : CG/JU
 Sample : N2471-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010083.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 3.48 ng/ul m

response 24990

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.88#
41.00	23.20	50.82#
42.00	18.00	31.52#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	170819	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	723336	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	501934	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1082903	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	914926	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	771631	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17962	4.585	ng/uL	0.00
4) Pyridine-d5	3.793	84	41280	3.746	ng/ul	0.00
7) Phenol-d5	7.093	99	87117	5.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	287626	30.851	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	261780	22.750	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	170694	13.260	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	177339	30.825	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	184701	29.397	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	299835	24.513	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	365893m	20.863	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1301269	33.328	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1400626	29.640	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	24990m	3.477	ng/ul	0.00
60) Fluorene-d10	15.563	176	1098366	32.702	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	187580	27.395	ng/ul	0.00
73) Anthracene-d10	17.422	188	1818120	34.704	ng/ul	0.00
81) Pyrene-d10	19.651	212	2065816	39.703	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1457009	36.017	ng/ul	0.00

Target Compounds Qvalue

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

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