

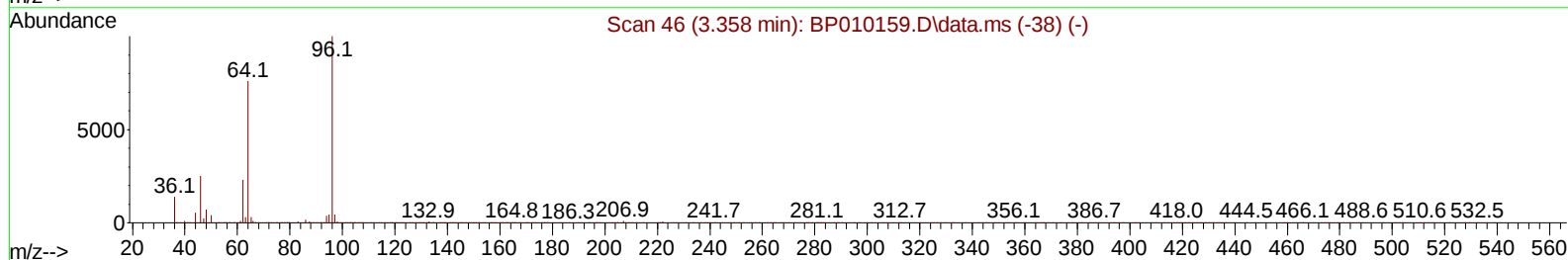
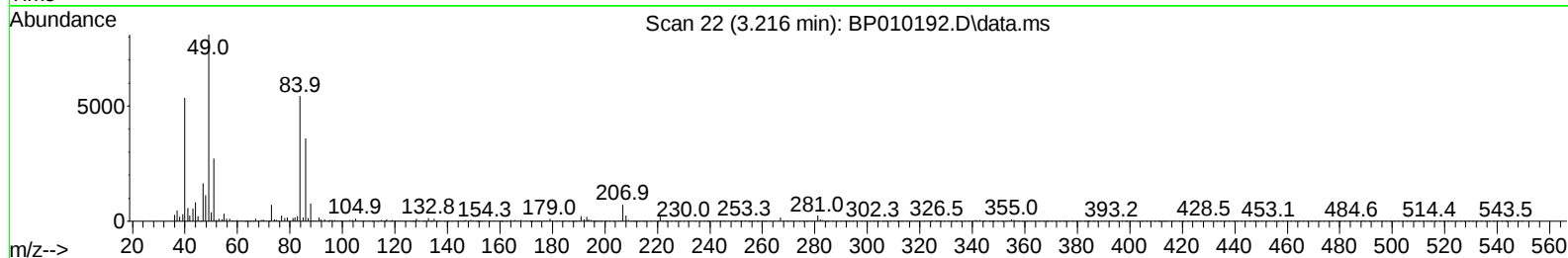
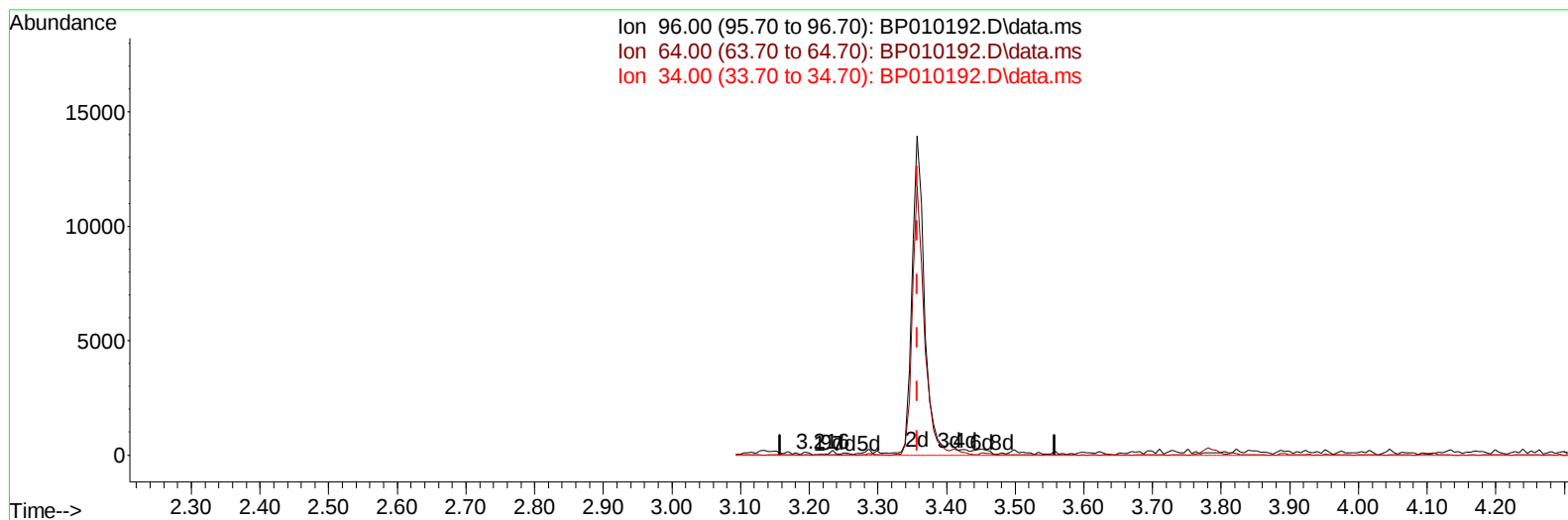
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010192.D
 Acq On : 03 May 2022 02:31
 Operator : CG/JU
 Sample : N2615-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0B1

Manual IntegrationsAPPROVED

Quant Time: May 03 03:24:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/03/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010192.D\data.ms

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

3.216min (-0.142) 0.01 ng/uL

response 22

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.60
34.00	0.00	0.00
0.00	0.00	0.00

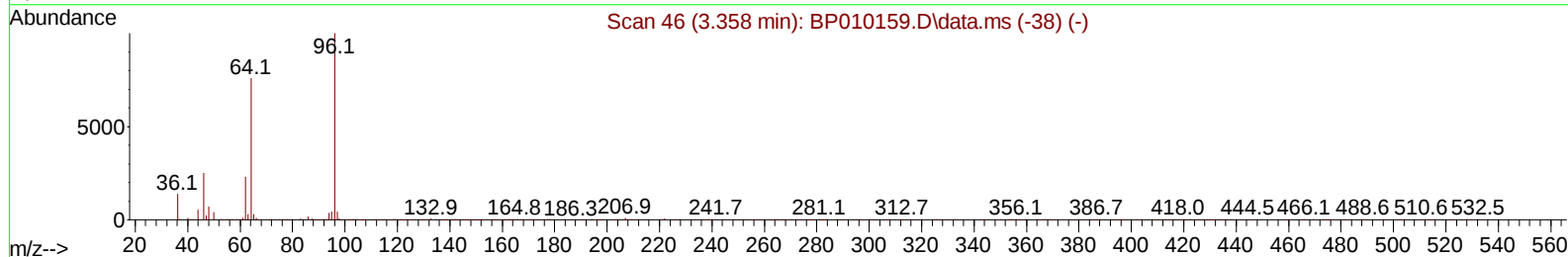
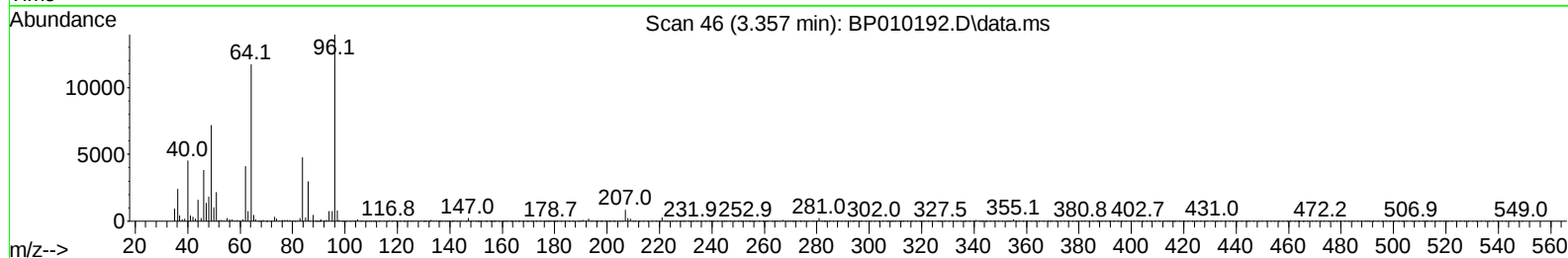
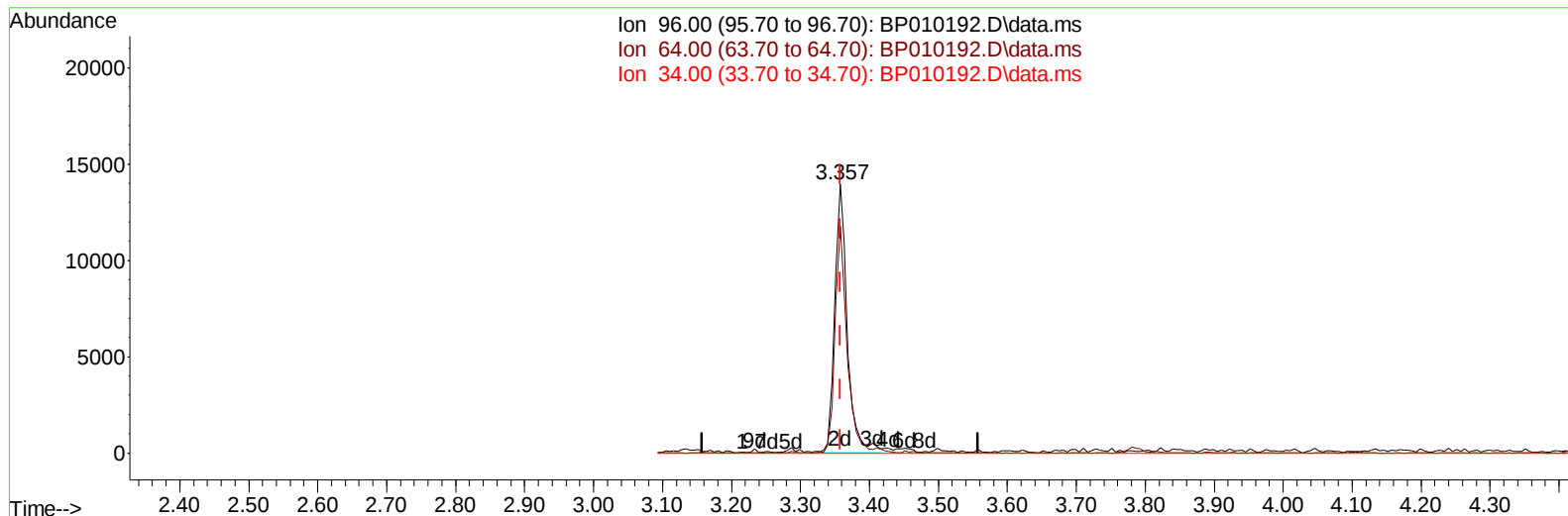
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010192.D
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Instrument :
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TIC: BP010192.D\data.ms

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

3.357min (-0.000) 5.62 ng/uL m

response 17388

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.25#
34.00	0.00	0.00
0.00	0.00	0.00

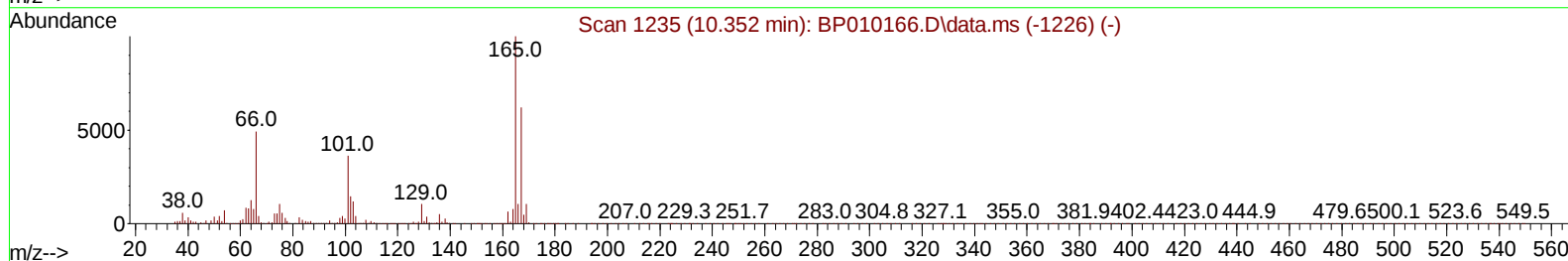
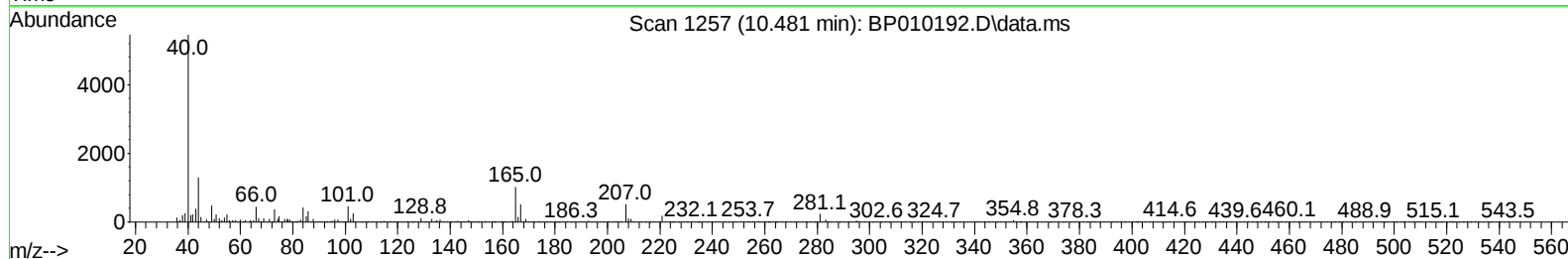
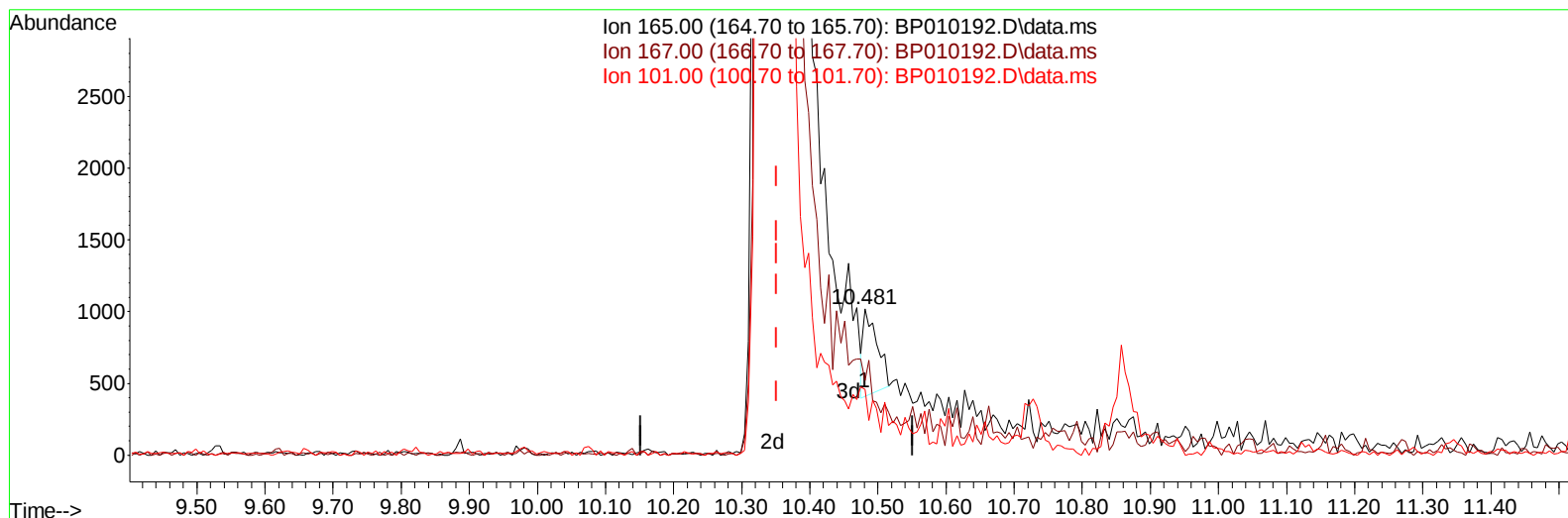
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TIC: BP010192.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.481min (+ 0.129) 0.09 ng/u1

response 852

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	51.13
101.00	49.80	44.94
0.00	0.00	0.00

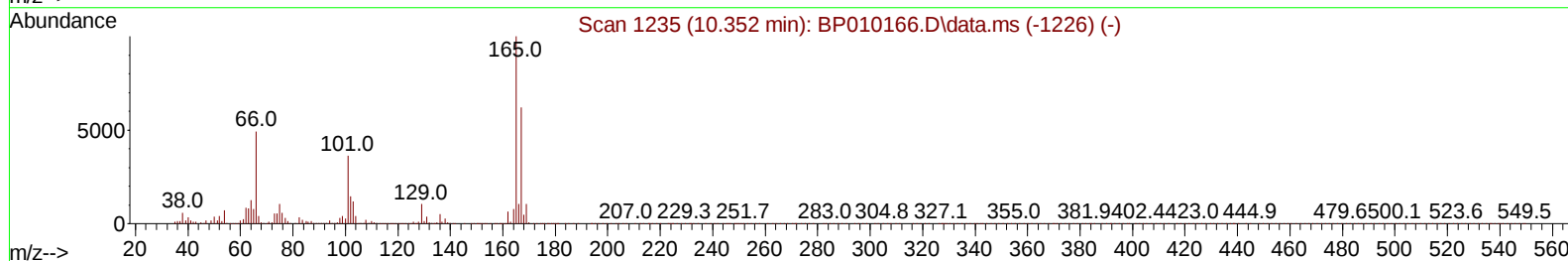
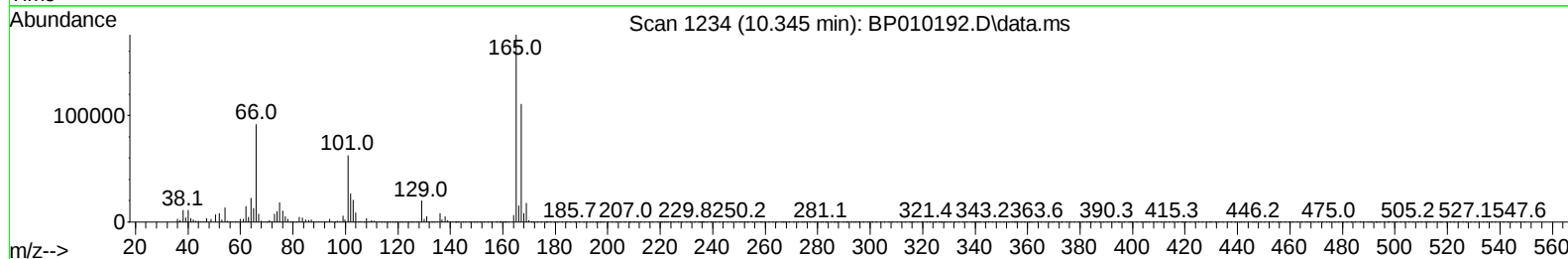
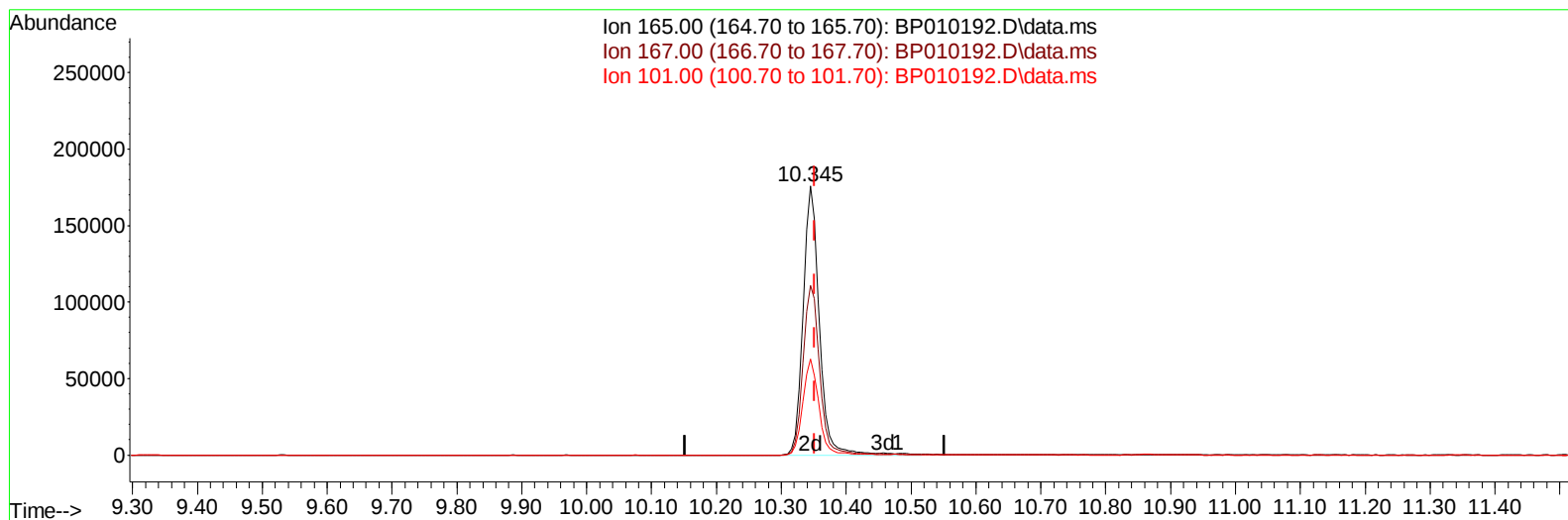
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Data File : BP010192.D
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Operator : CG/JU
Sample : N2615-15
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0B1

Manual IntegrationsAPPROVED

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TIC: BP010192.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.345min (-0.006) 32.84 ng/ul m

response 307143

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.14
101.00	49.80	35.70#
0.00	0.00	0.00

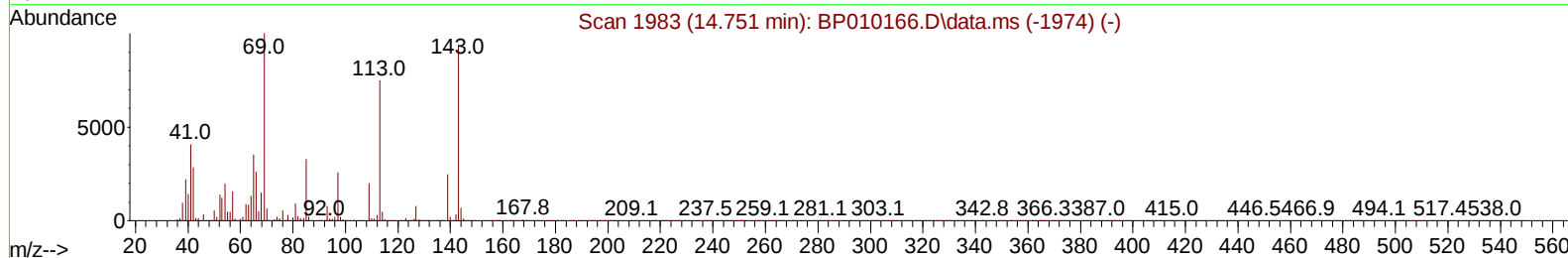
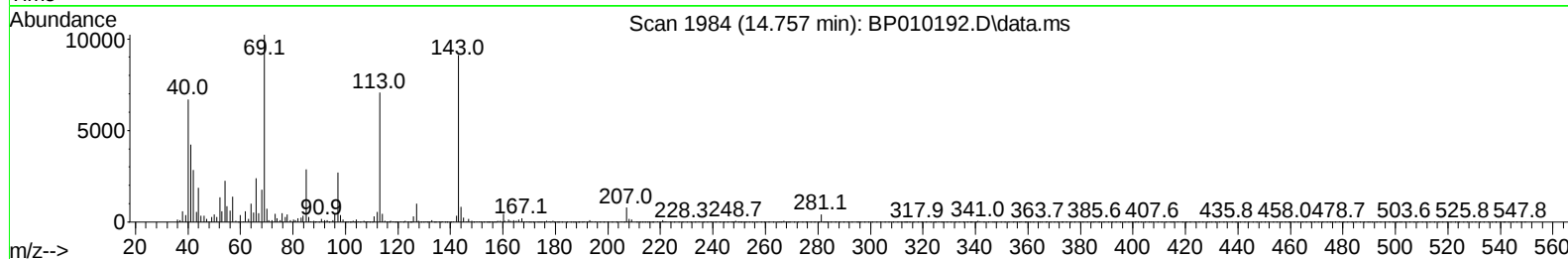
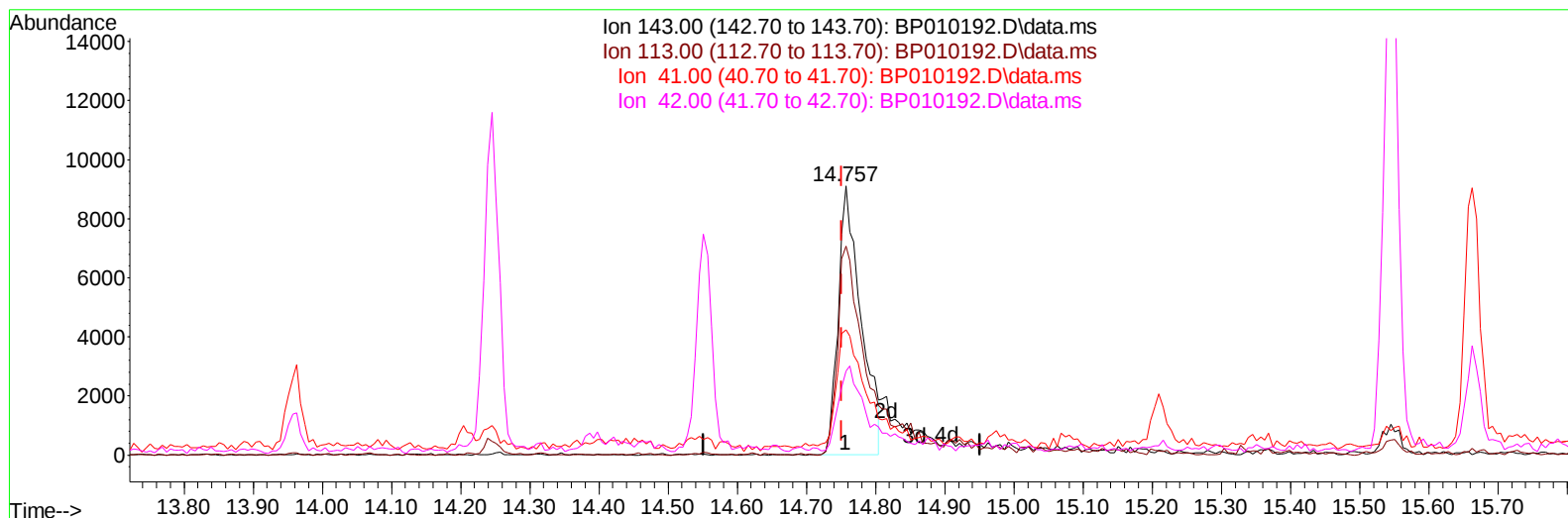
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010192.D
Acq On : 03 May 2022 02:31
Operator : CG/JU
Sample : N2615-15
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH0B1

Manual IntegrationsAPPROVED

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TIC: BP010192.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 4.41 ng/ul

response 20790

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.58#
41.00	23.20	46.52#
42.00	18.00	31.16#

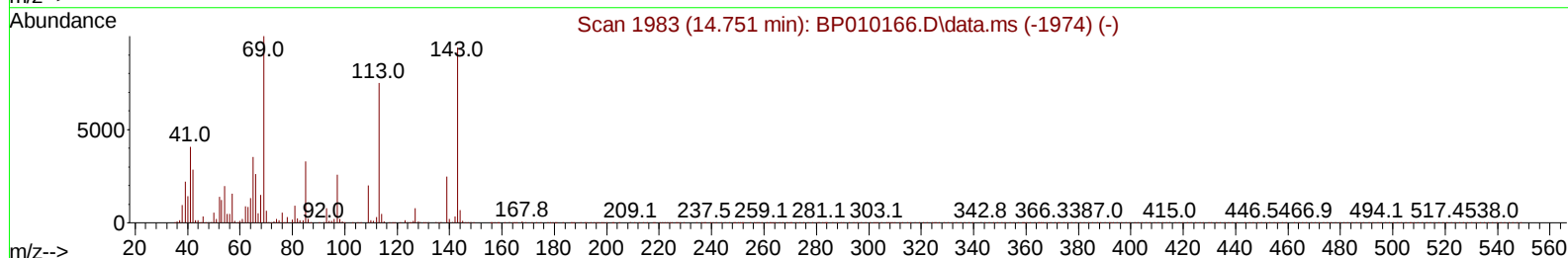
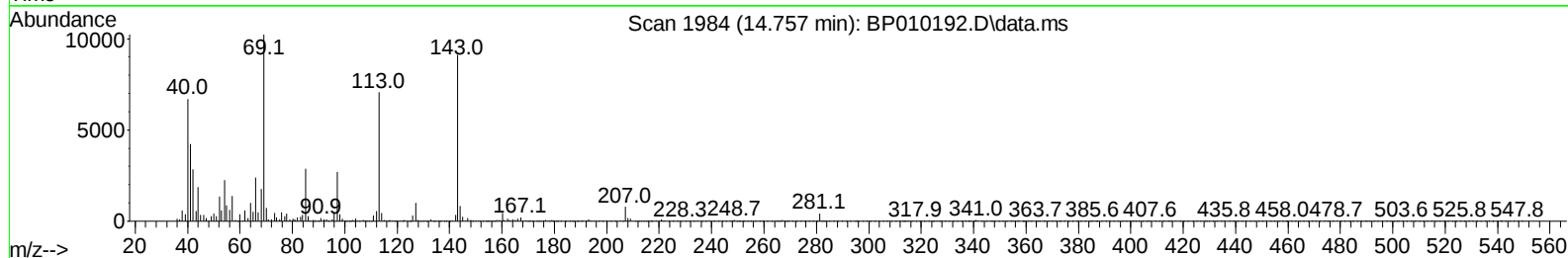
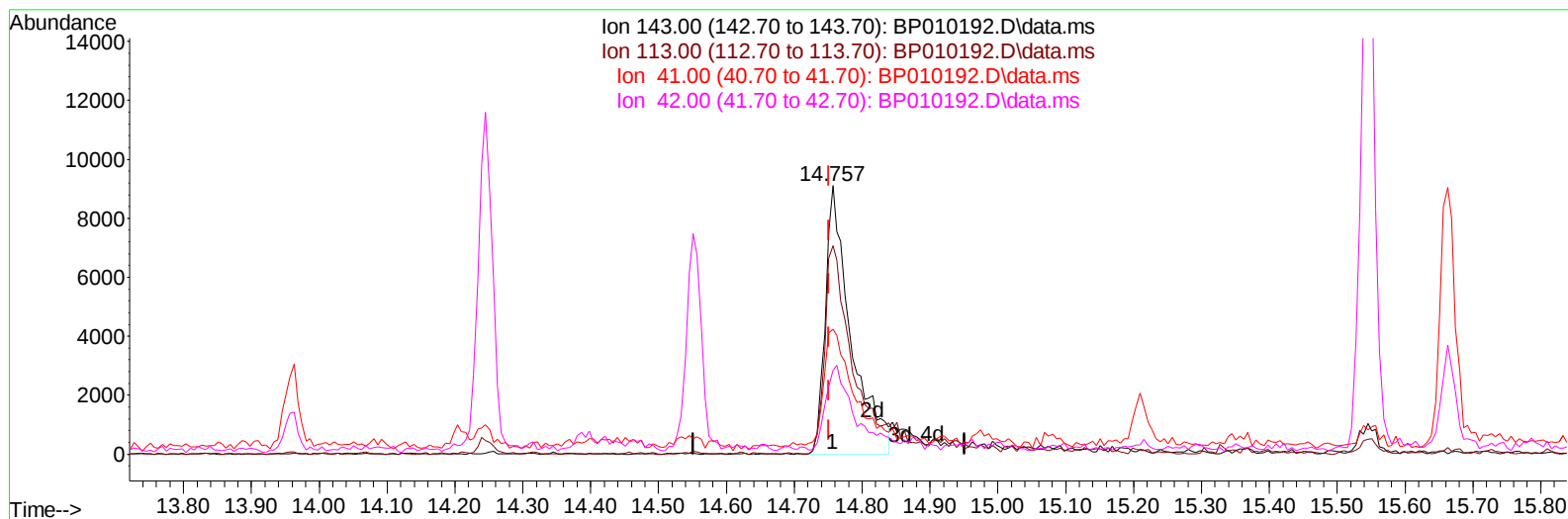
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010192.D
 Acq On : 03 May 2022 02:31
 Operator : CG/JU
 Sample : N2615-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0B1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/03/2022
 Supervised By :Yogesh Patel 05/05/2022

Quant Time: May 03 03:24:37 2022
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 Response via : Initial Calibration



TIC: BP010192.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 5.04 ng/ul m

response 23766

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.58#
41.00	23.20	46.52#
42.00	18.00	31.16#

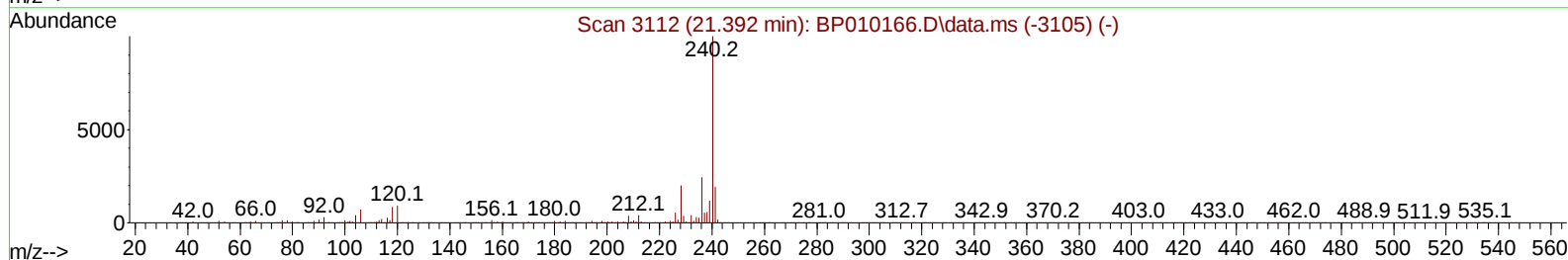
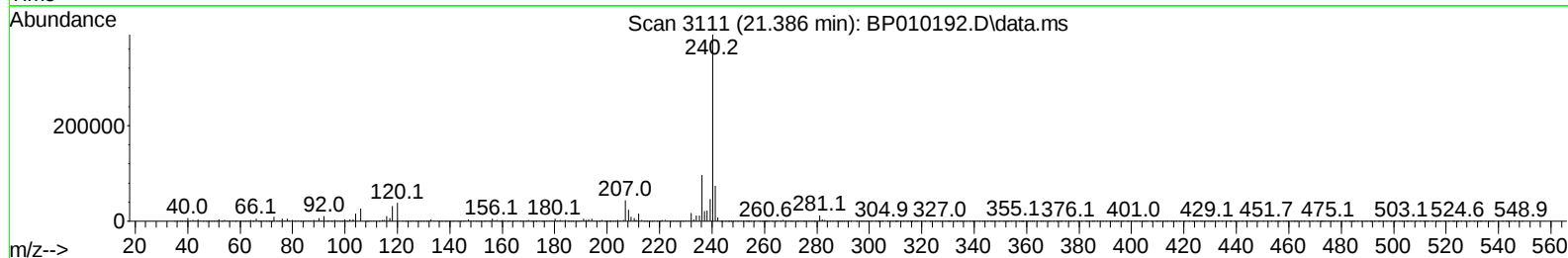
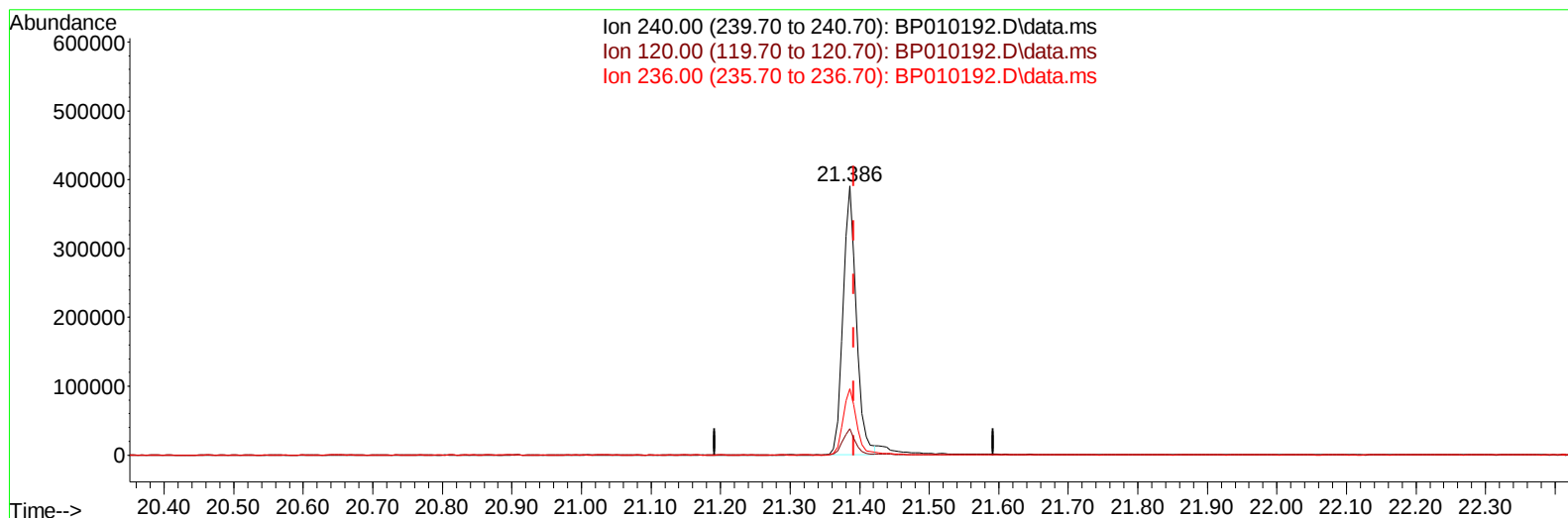
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 Misc :
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Instrument :
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TIC: BP010192.D\data.ms

(79) Chrysene-d12 (I)

21.386min (-0.006) 20.00 ng/u1

response 523088

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.71
236.00	31.60	24.72#
0.00	0.00	0.00

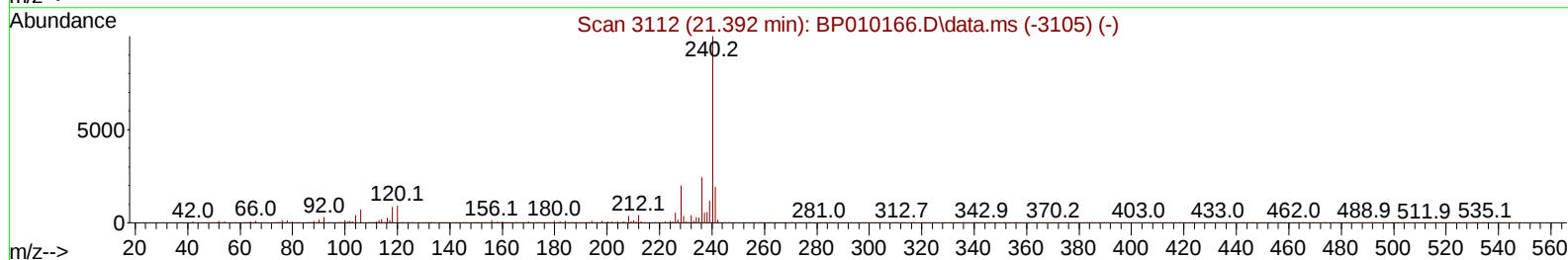
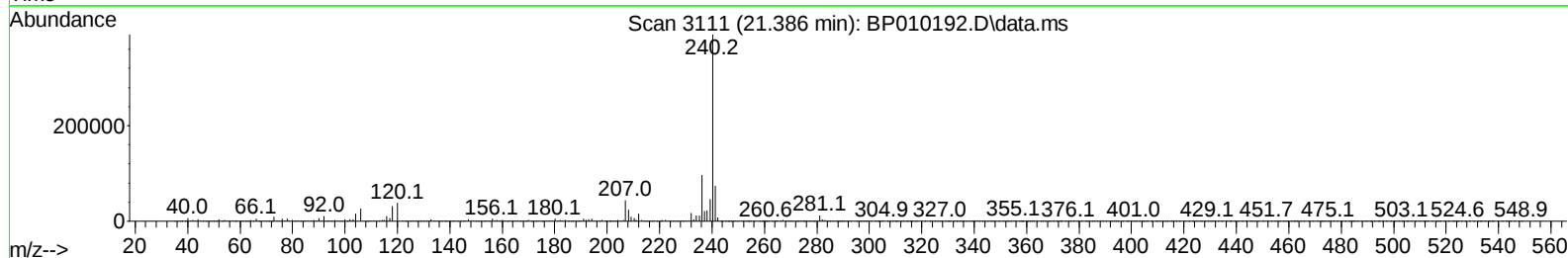
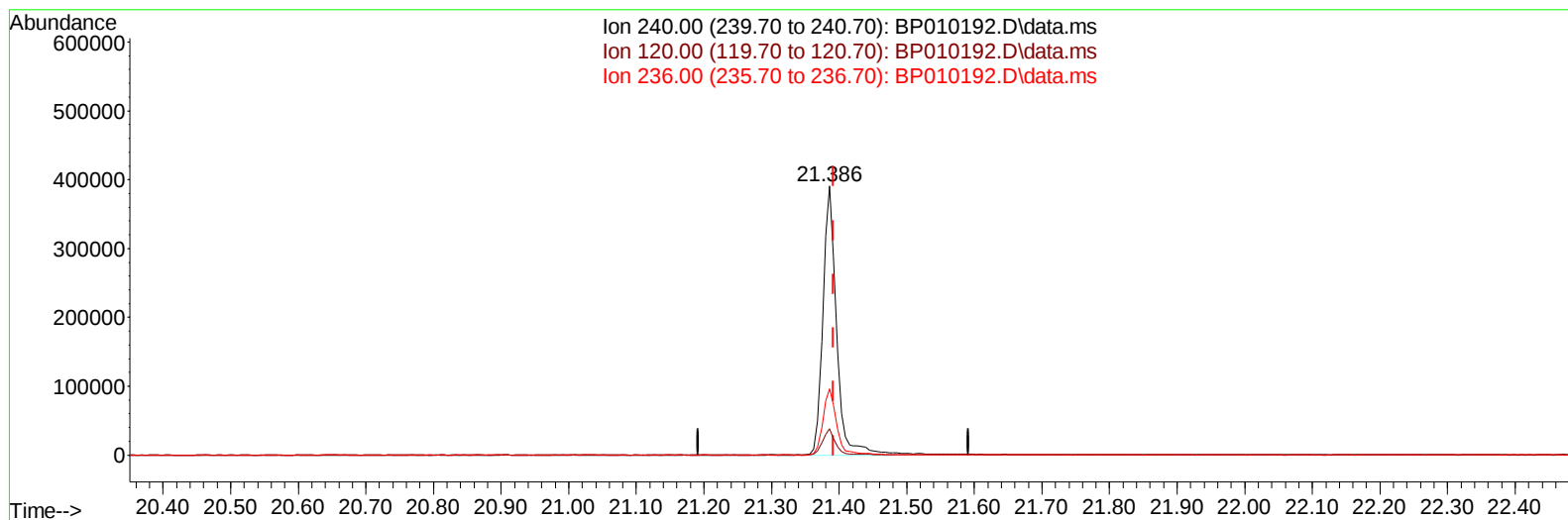
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
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 Operator : CG/JU
 Sample : N2615-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP010192.D\data.ms

(79) Chrysene-d12 (I)

21.386min (-0.006) 20.00 ng/ul m

response 548523

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.71
236.00	31.60	24.72#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	134986	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.728	136	553056	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	329451	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	618565	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	548523m	20.000	ng/ul	0.00
88) Perylene-d12	23.827	264	539082	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	17388m	5.617	ng/uL	0.00
4) Pyridine-d5	3.781	84	98521	11.313	ng/ul	0.00
7) Phenol-d5	7.087	99	88086	7.174	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.245	67	307998	41.805	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	270412	29.738	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	176126	17.313	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	187939	42.726	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	196003	40.800	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	307143m	32.842	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	388394	28.964	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1152525	44.972	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1291789	41.649	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	23766m	5.037	ng/ul	0.00
60) Fluorene-d10	15.545	176	964541	43.753	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	152758	39.056	ng/ul	0.00
73) Anthracene-d10	17.404	188	1487778	49.716	ng/ul	0.00
81) Pyrene-d10	19.633	212	1678989	53.823	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1486677	52.604	ng/ul	0.00

Target Compounds Qvalue

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

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