

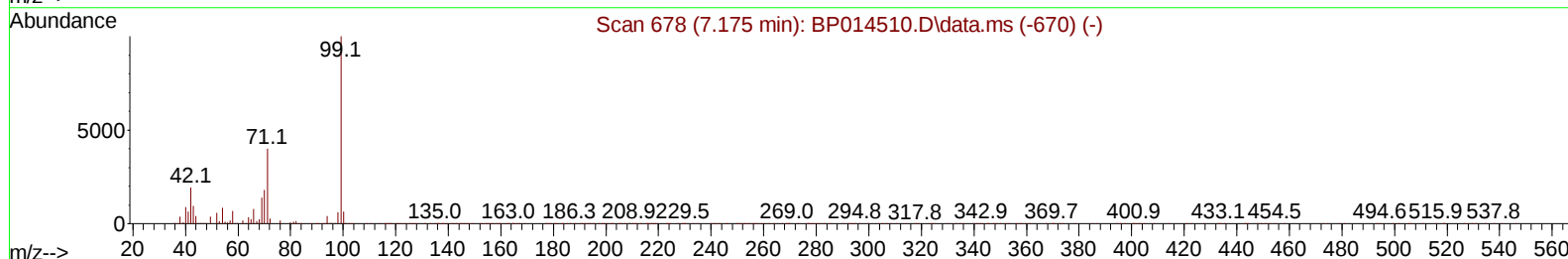
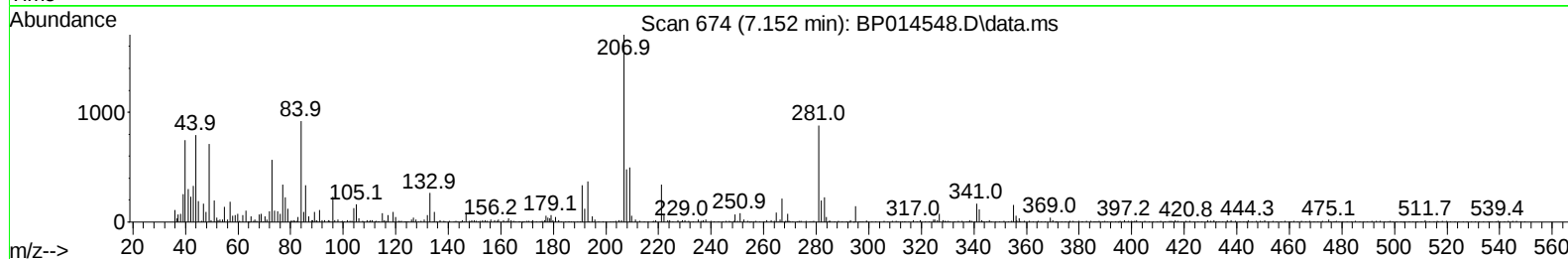
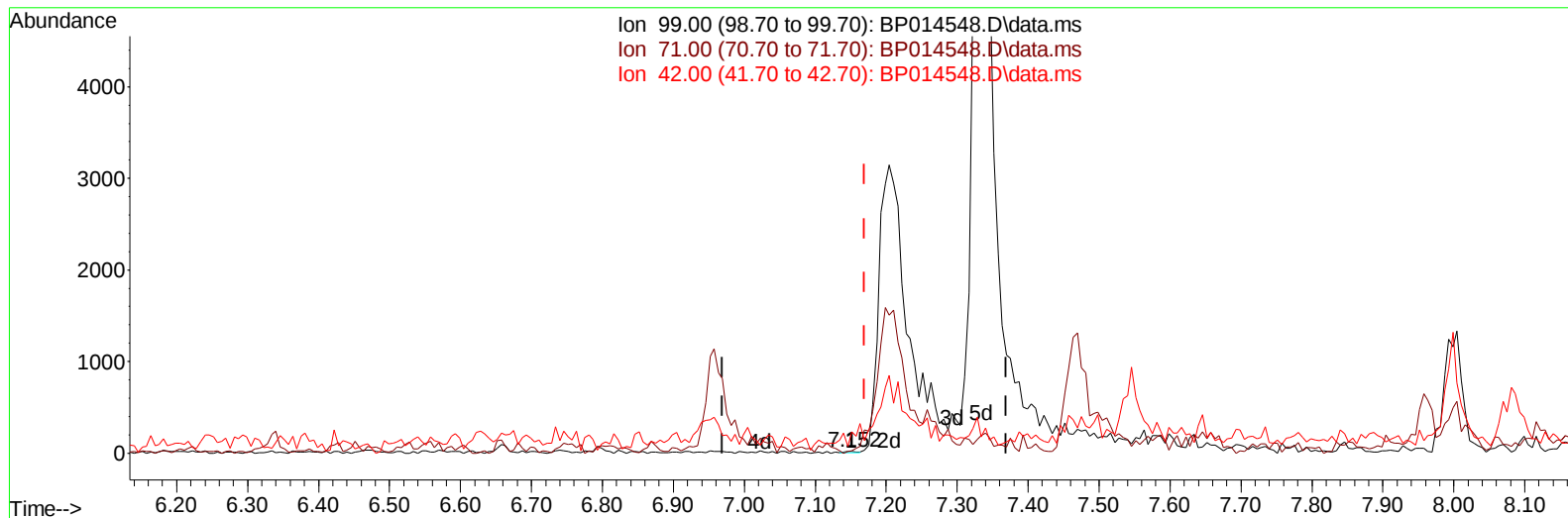
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014548.D
Acq On : 05 Apr 2023 03:06
Operator : CG/JU
Sample : 02123-02DL 10X
Misc :
ALS Vial : 72 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9A2DL

Manual IntegrationsAPPROVED

Quant Time: Apr 05 03:51:05 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 05 02:40:56 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/05/2023
Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014548.D\data.ms

(7) Phenol-d5 (S)

7.152min (-0.018) 0.00 ng/u1

response 14

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	137.50#
42.00	18.90	1450.00#
0.00	0.00	0.00

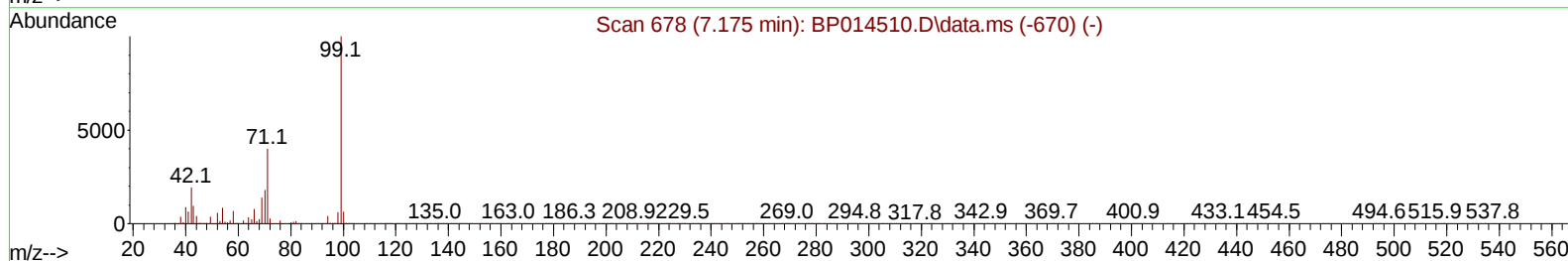
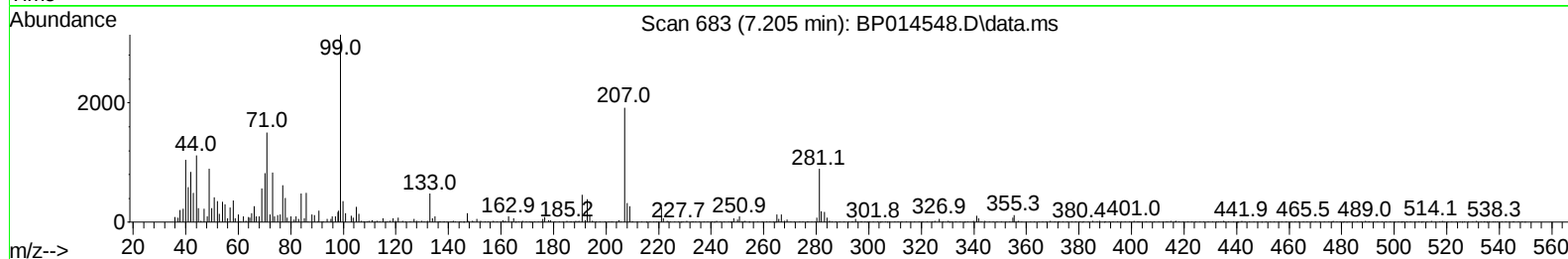
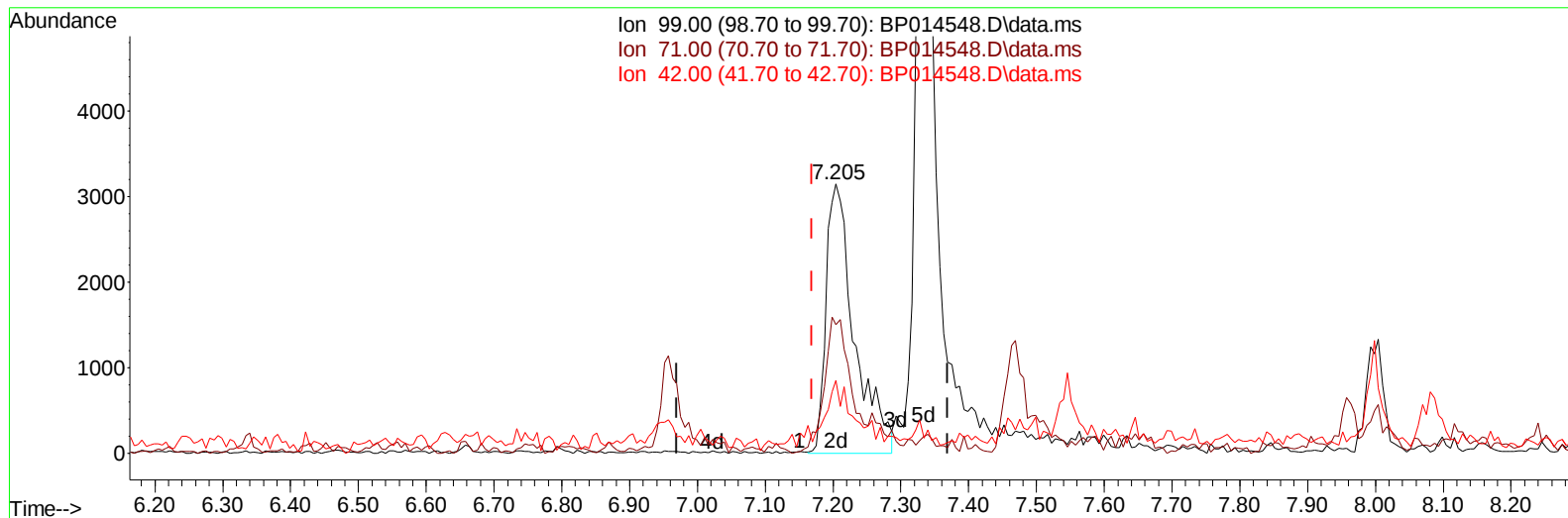
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(7) Phenol-d5 (S)

7.205min (+ 0.035) 0.91 ng/ul m

response 9128

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	47.84
42.00	18.90	26.92#
0.00	0.00	0.00

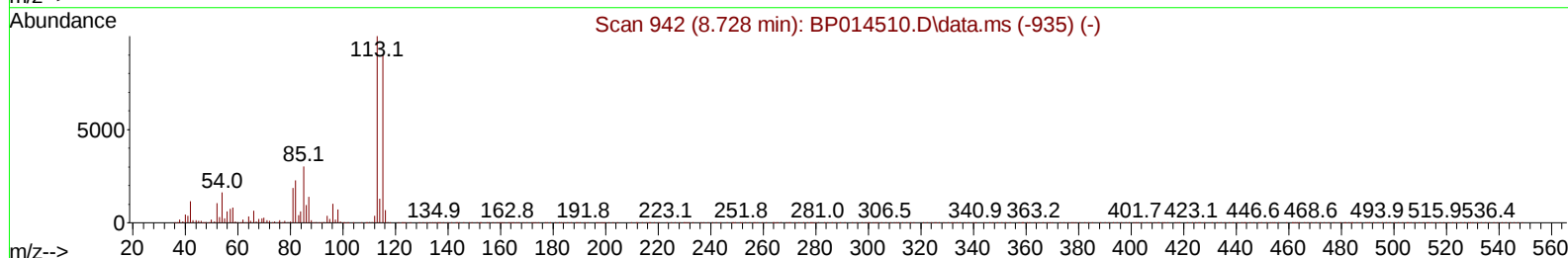
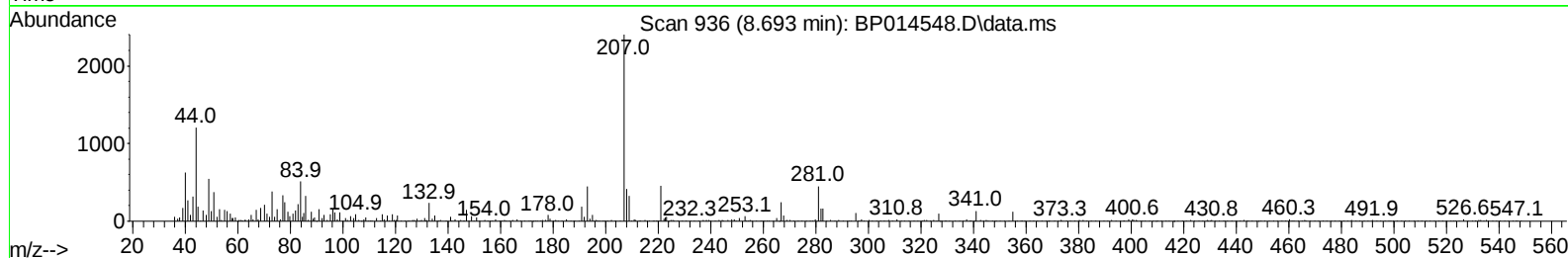
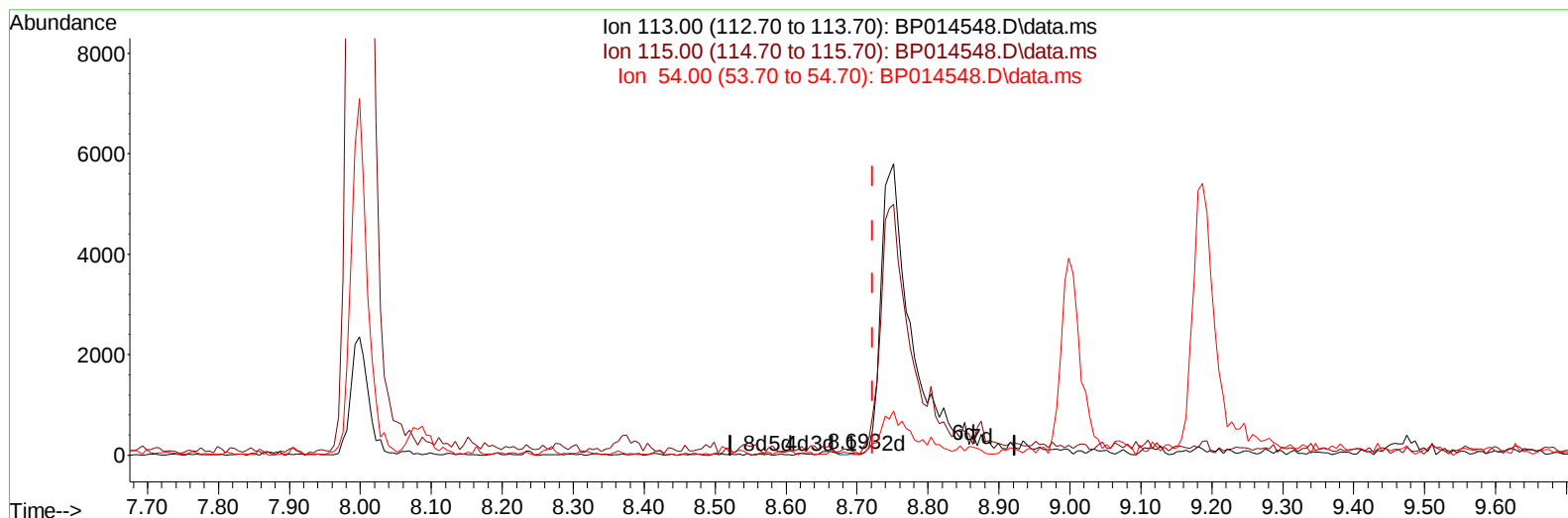
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TIC: BP014548.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.693min (-0.030) 0.00 ng/u1

response 25

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	89.90	223.08#
54.00	15.10	38.46#
0.00	0.00	0.00

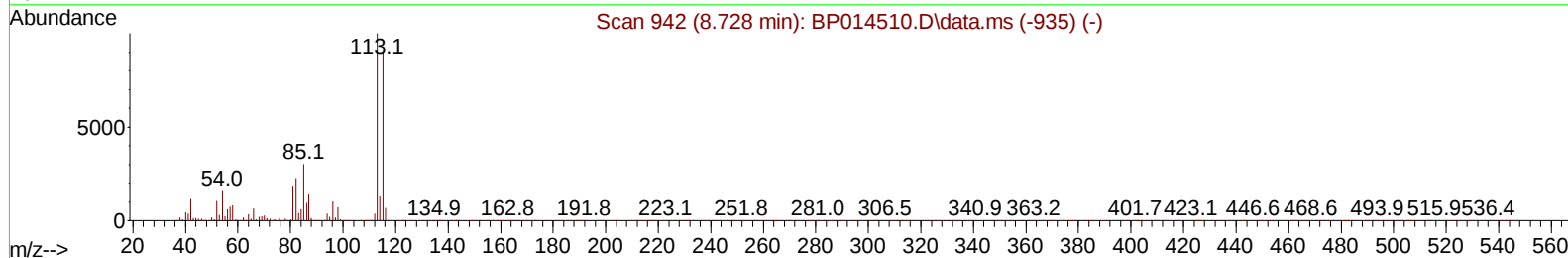
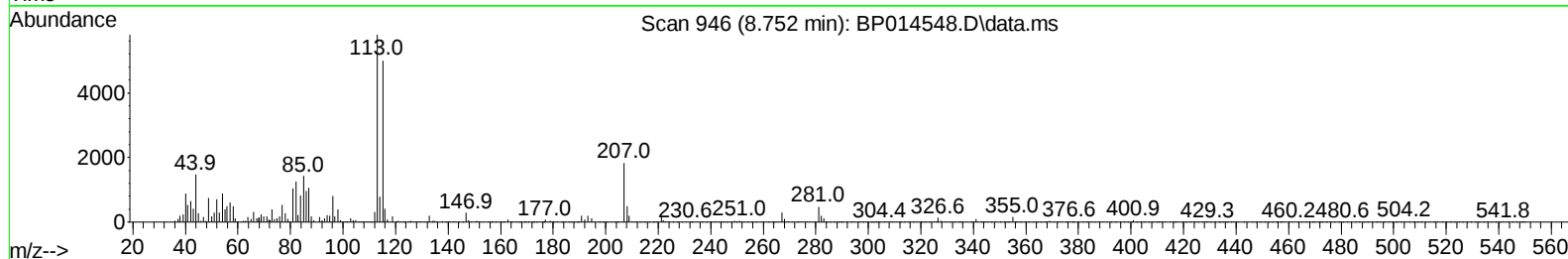
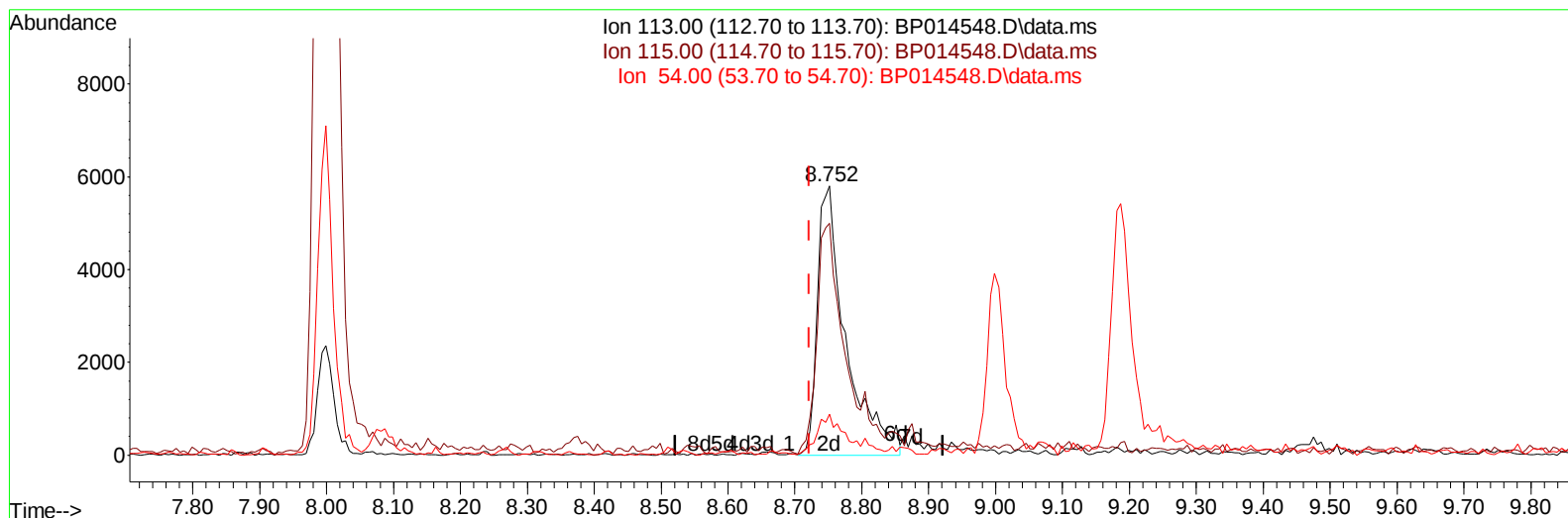
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TIC: BP014548.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.752min (+ 0.029) 2.11 ng/u1 m

response 17134

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	89.90	86.17
54.00	15.10	15.23
0.00	0.00	0.00

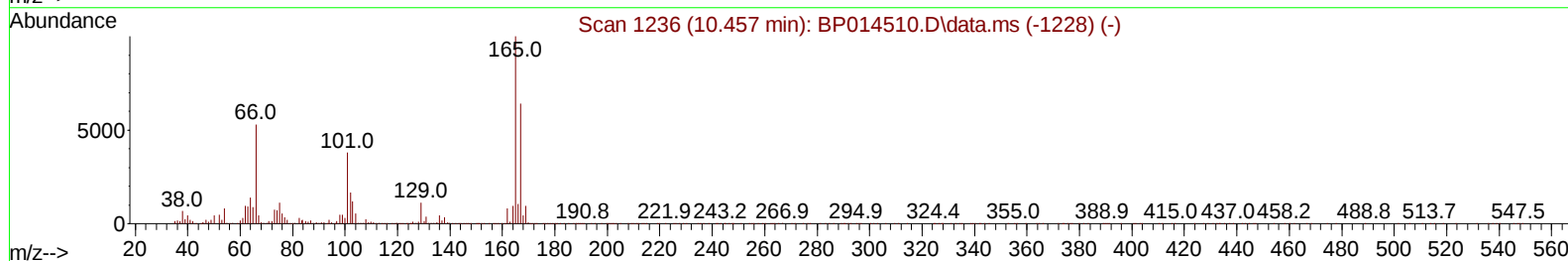
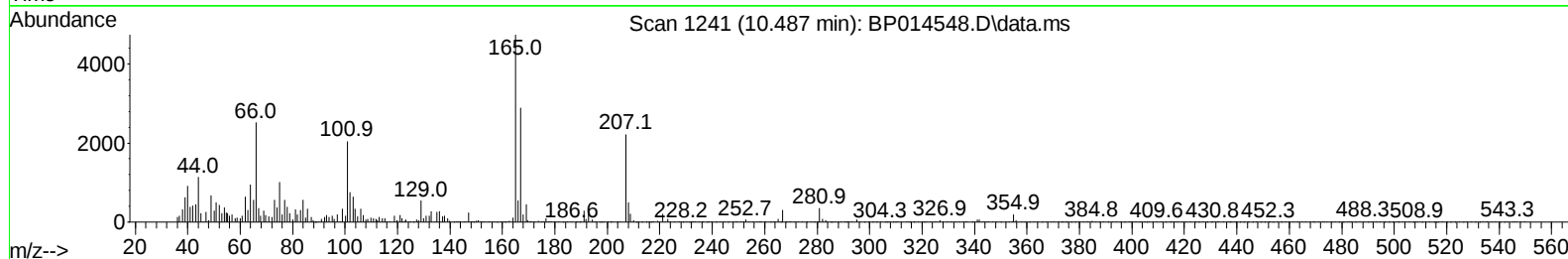
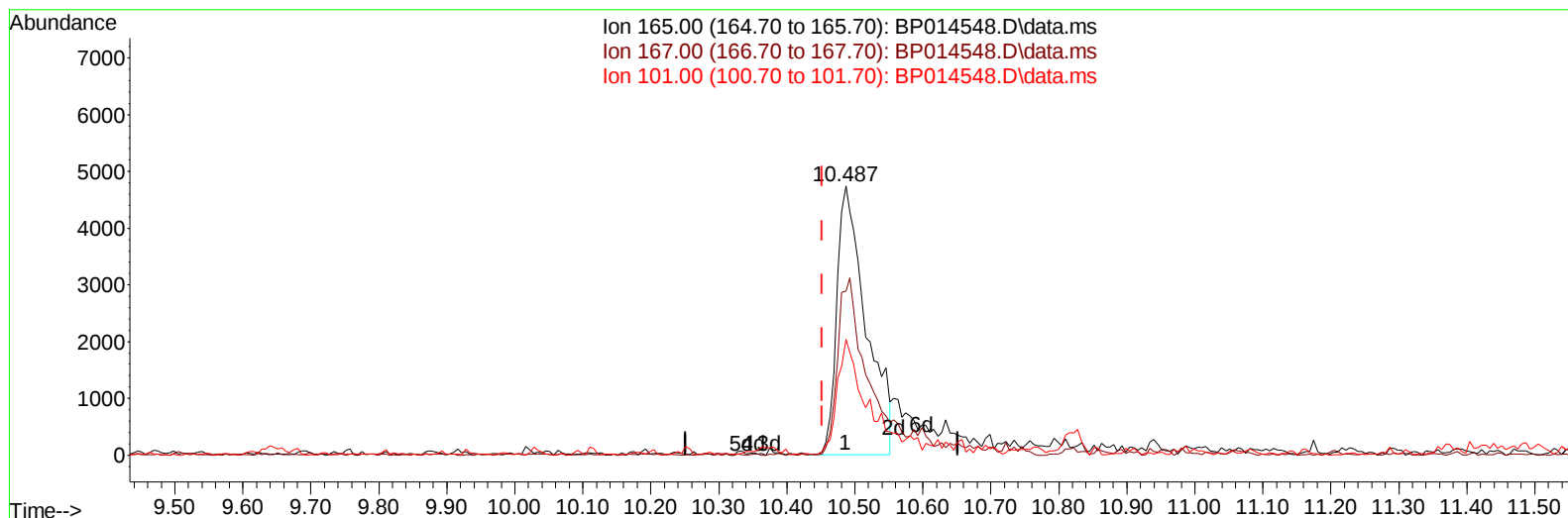
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Data File : BP014548.D
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Operator : CG/JU
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Misc :
ALS Vial : 72 Sample Multiplier: 1

Instrument :
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TIC: BP014548.D\data.ms

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

10.487min (+ 0.035) 1.86 ng/u1

response 14163

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.80	60.88
101.00	39.20	43.10
0.00	0.00	0.00

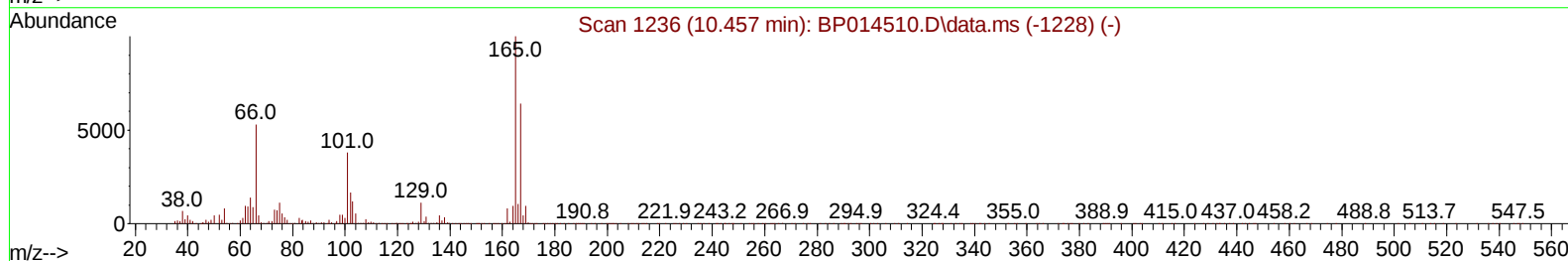
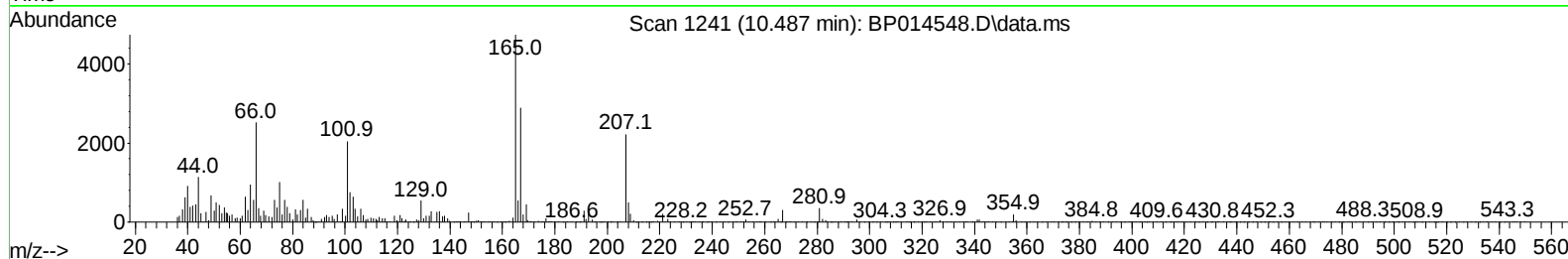
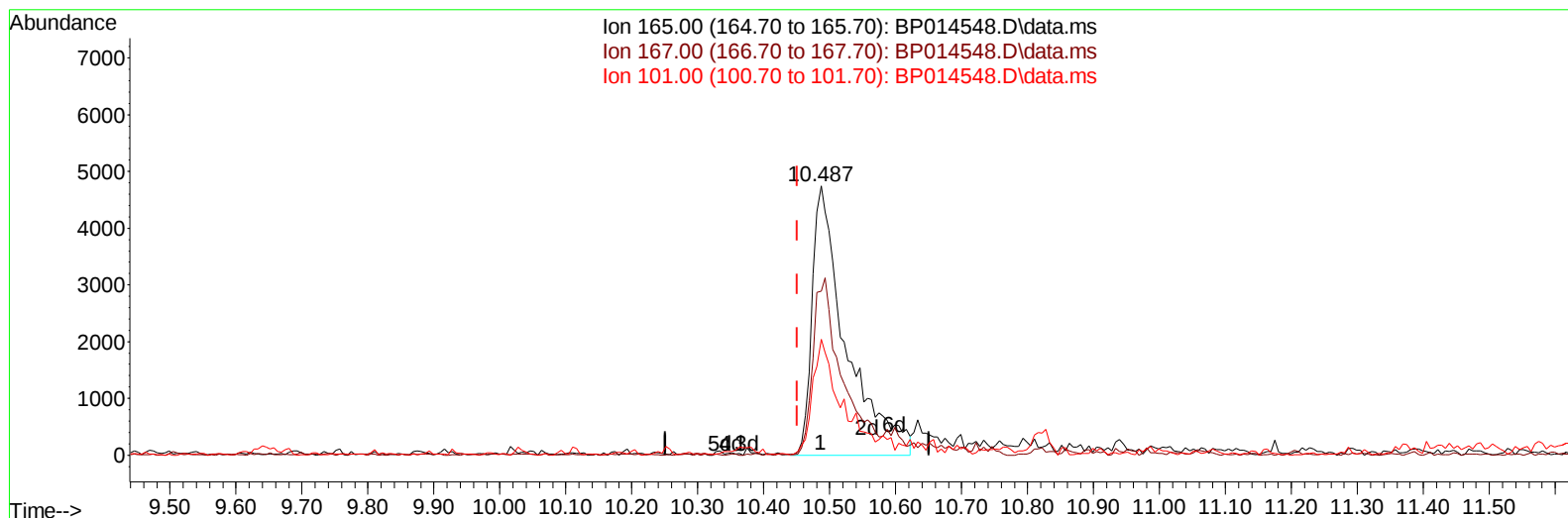
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014548.D
Acq On : 05 Apr 2023 03:06
Operator : CG/JU
Sample : 02123-02DL 10X
Misc :
ALS Vial : 72 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP014548.D\data.ms

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

10.487min (+ 0.035) 2.21 ng/ul m

response 16832

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.80	60.88
101.00	39.20	43.10
0.00	0.00	0.00

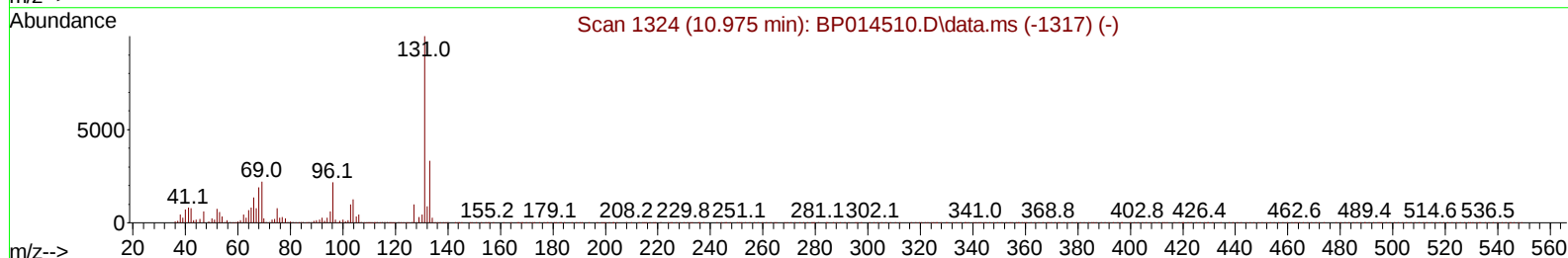
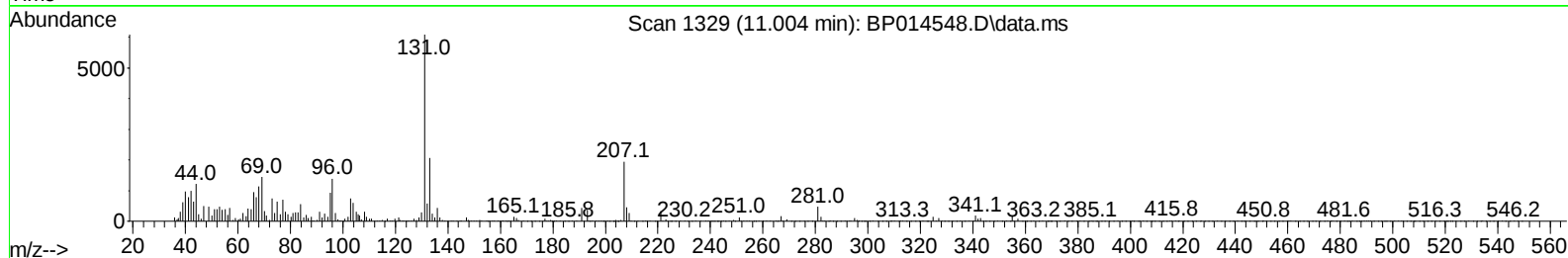
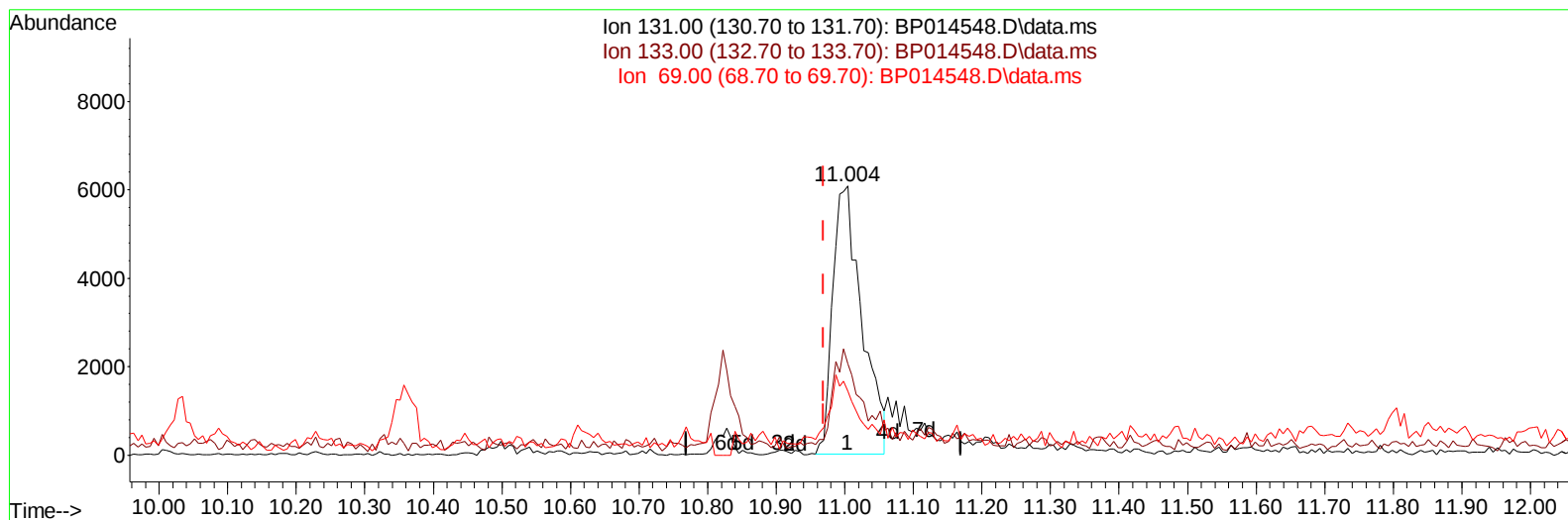
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TIC: BP014548.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.004min (+ 0.035) 1.78 ng/u1

response 17898

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.91
69.00	21.30	23.69
0.00	0.00	0.00

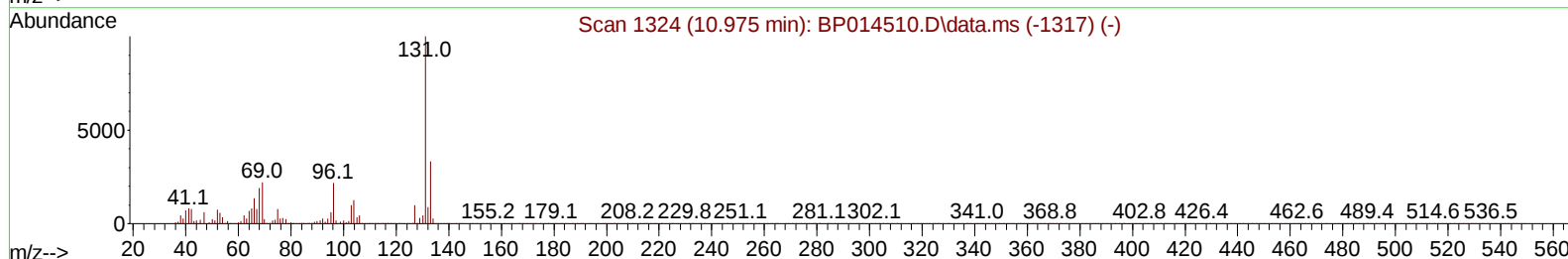
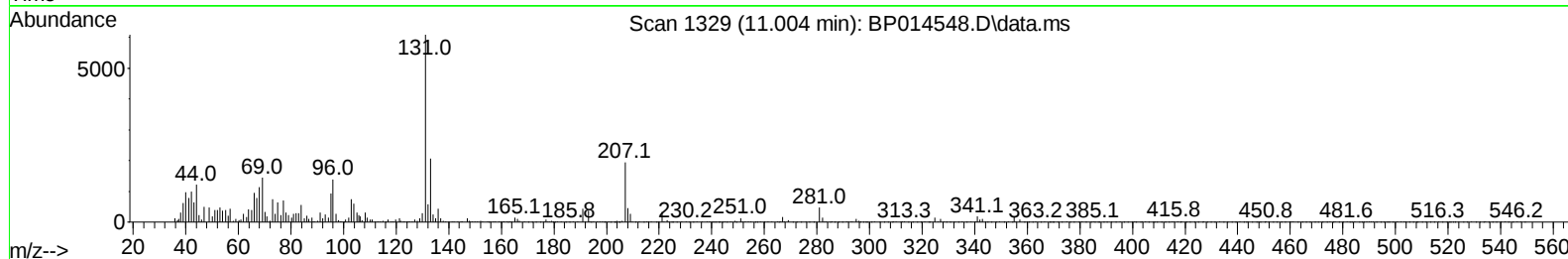
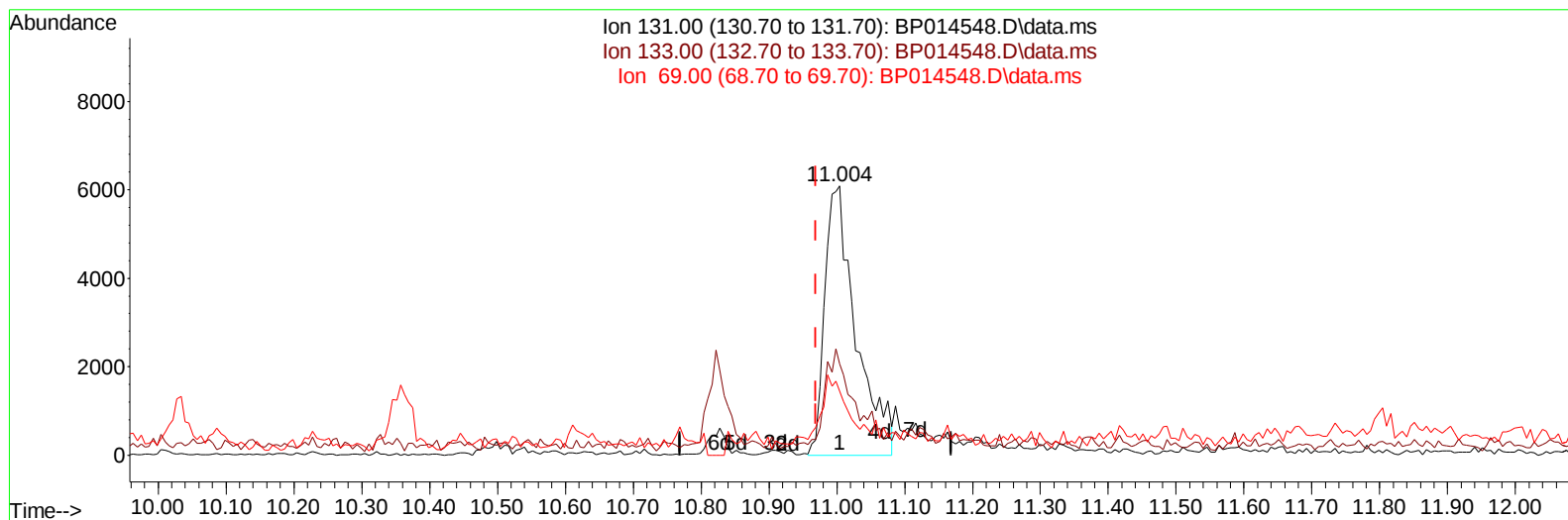
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Misc :
ALS Vial : 72 Sample Multiplier: 1

Instrument :
BNA_P
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TIC: BP014548.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.004min (+ 0.035) 1.93 ng/ul m

response 19429

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.91
69.00	21.30	23.69
0.00	0.00	0.00

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 Misc :
 ALS Vial : 72 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.999	152	109167	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	452921	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	286128	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	667765	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	718400	20.000	ng/ul	0.00
88) Perylene-d12	24.010	264	835513	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	1809	0.643	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.205	99	9128m	0.907	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.334	67	14828	2.295	ng/ul	0.01
11) 2-Chlorophenol-d4	7.540	132	13898	1.910	ng/ul	0.01
15) 4-Methylphenol-d8	8.752	113	17134m	2.110	ng/ul	0.03
21) Nitrobenzene-d5	9.187	128	7876	2.154	ng/ul	0.01
24) 2-Nitrophenol-d4	9.916	143	7416	1.771	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.487	165	16832m	2.207	ng/ul	0.04
31) 4-Chloroaniline-d4	11.004	131	19429m	1.934	ng/ul	0.04
46) Dimethylphthalate-d6	14.063	166	67834	2.924	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	71522	2.778	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.651	176	60207	3.112	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.516	188	101846	3.151	ng/ul	0.00
81) Pyrene-d10	19.745	212	134130	2.780	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	135875	3.061	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.417	88	80873	26.142	ng/uL	98

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

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