

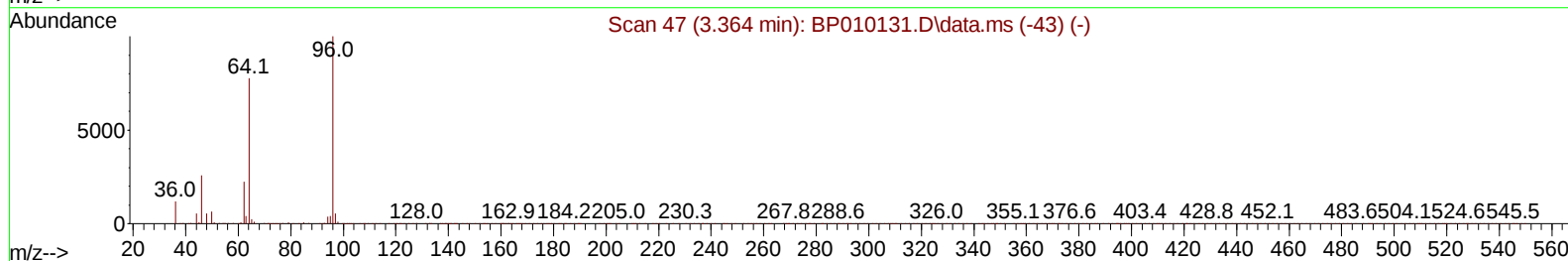
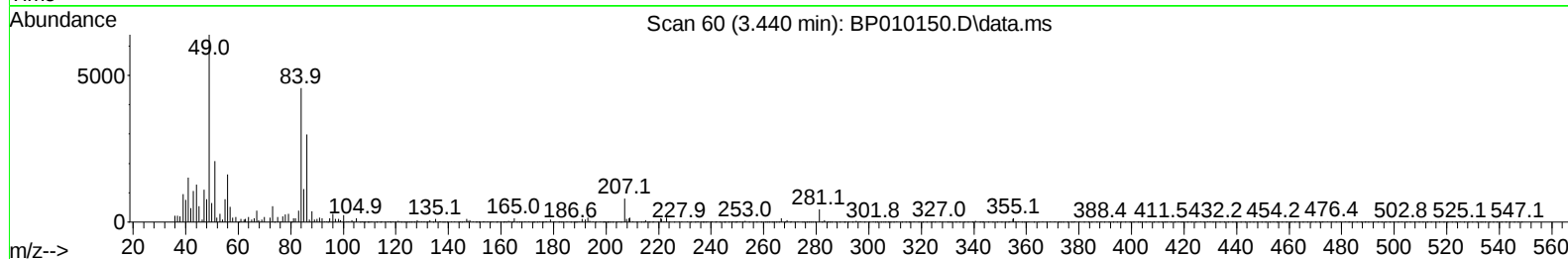
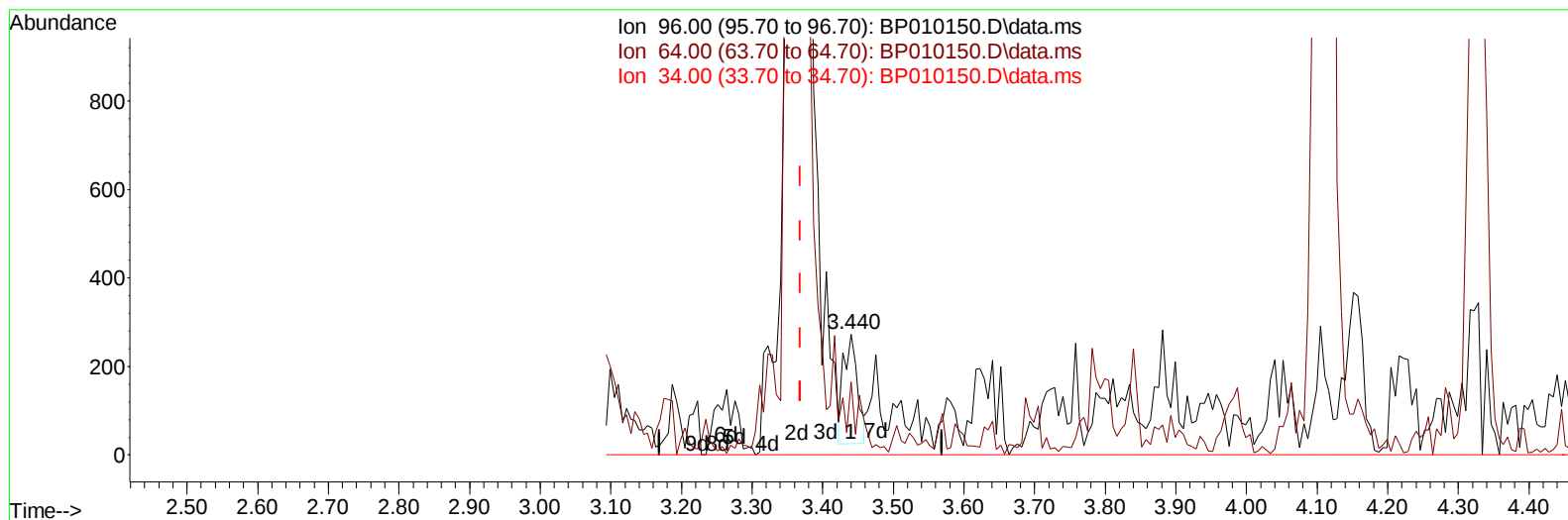
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010150.D
 Acq On : 29 Apr 2022 14:28
 Operator : CG/JU
 Sample : N2587-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXET7

Manual IntegrationsAPPROVED

Quant Time: Apr 30 00:30:19 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
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Reviewed By :Jagrut Upadhyay 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010150.D\data.ms

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

3.440min (+ 0.071) 0.10 ng/uL

response 331

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	60.81
34.00	0.00	0.00
0.00	0.00	0.00

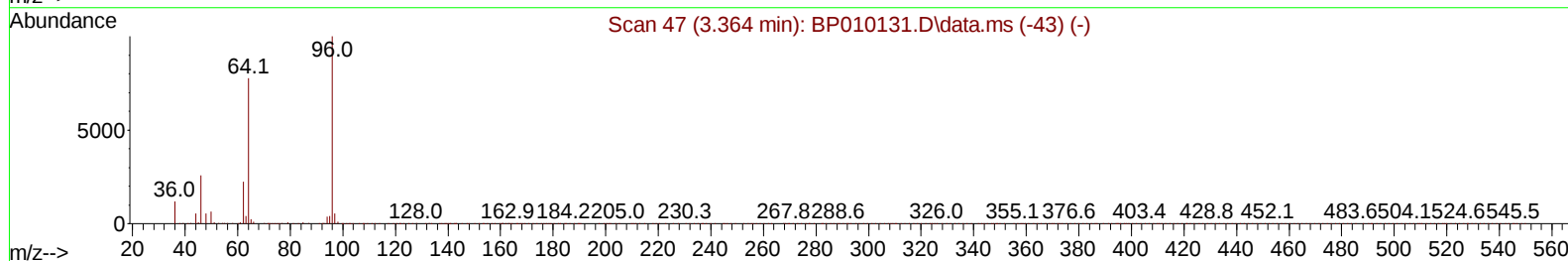
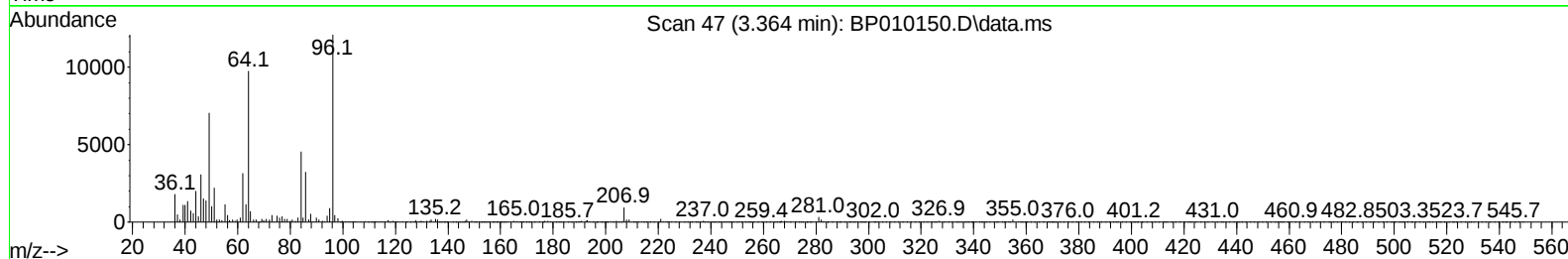
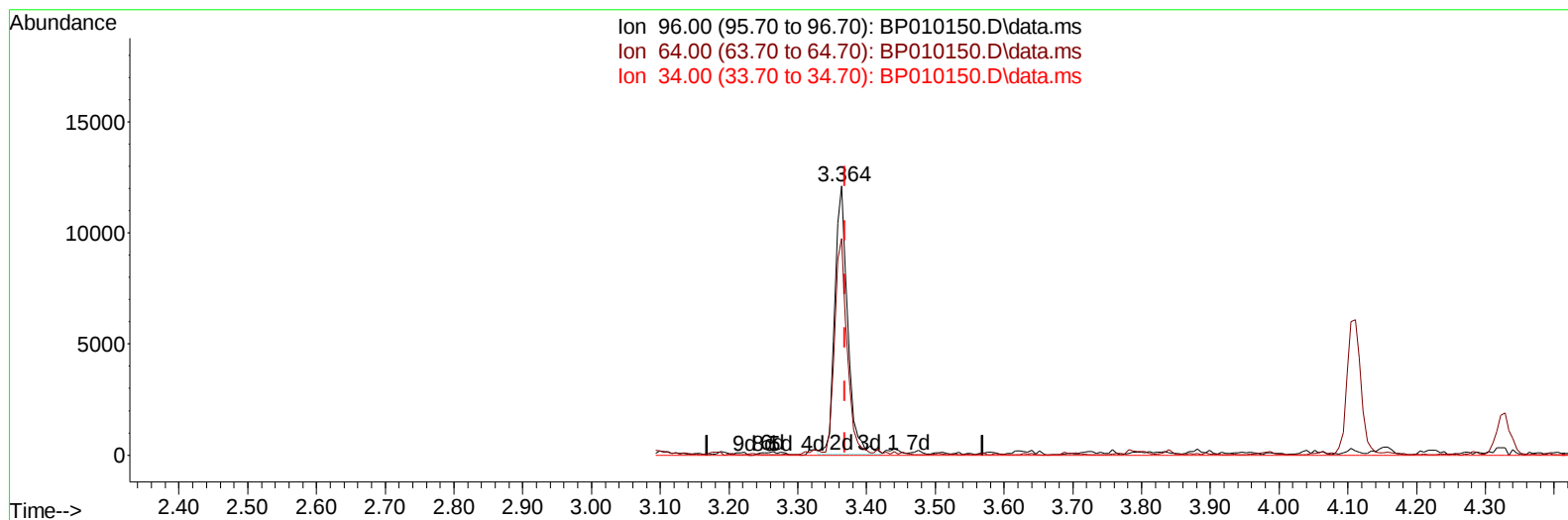
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Instrument :
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TIC: BP010150.D\data.ms

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

3.364min (-0.005) 4.74 ng/uL m

response 15823

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.60#
34.00	0.00	0.00
0.00	0.00	0.00

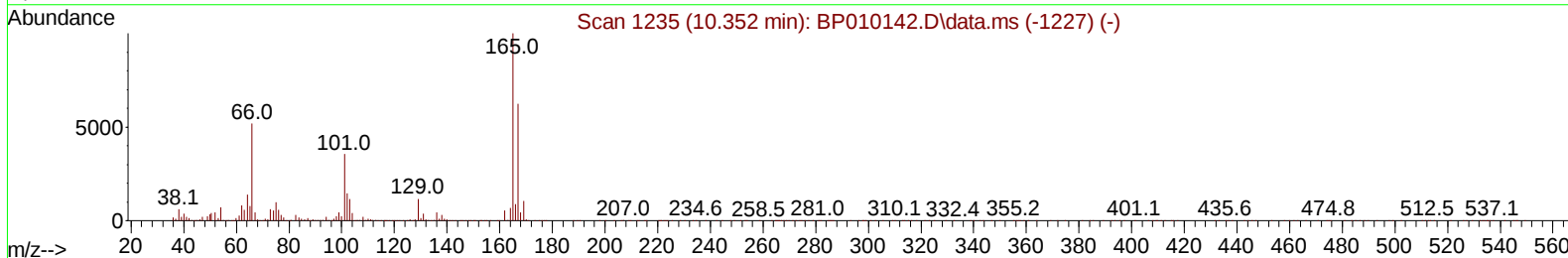
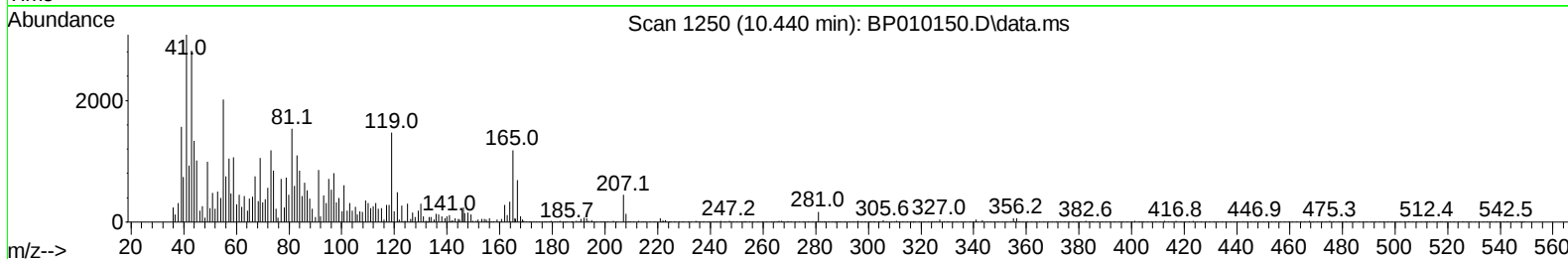
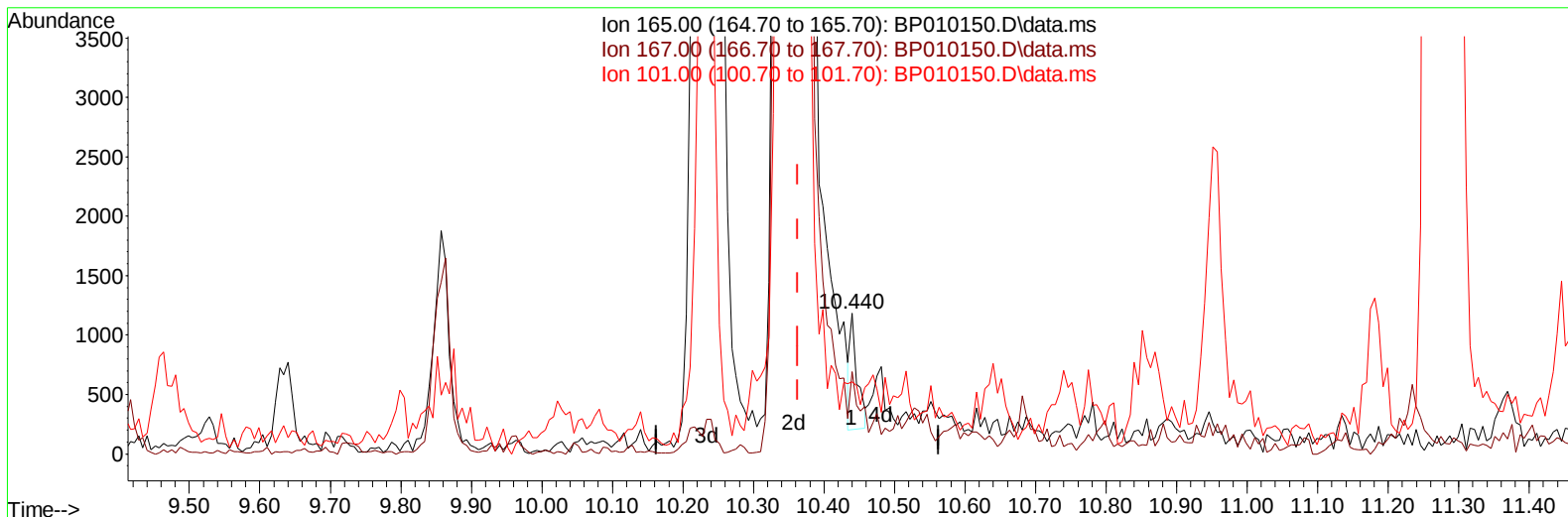
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TIC: BP010150.D\data.ms

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

10.440min (+ 0.077) 0.07 ng/u1

response 662

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	58.35
101.00	49.80	51.26
0.00	0.00	0.00

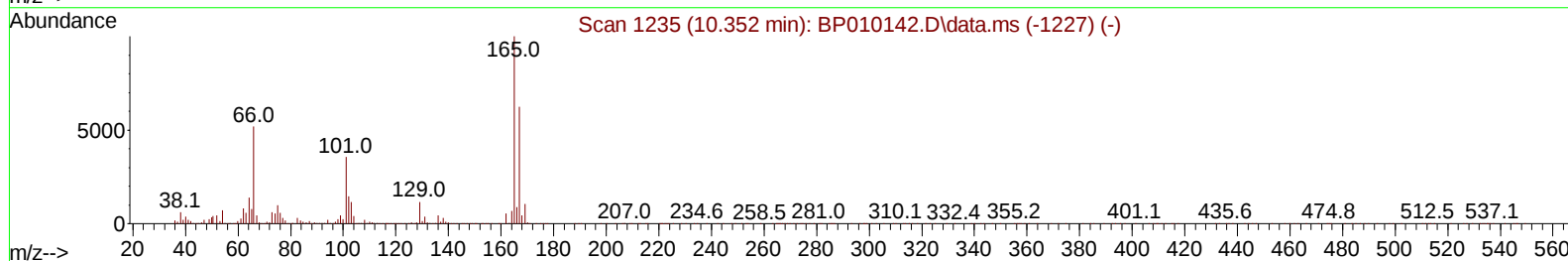
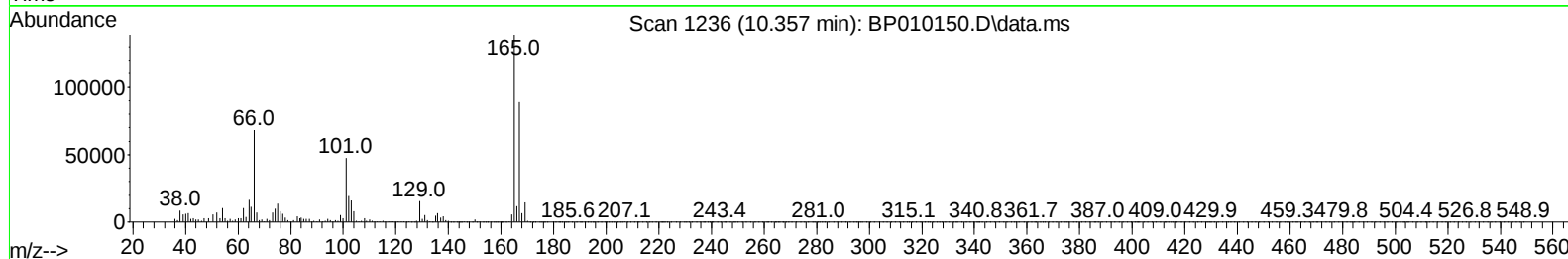
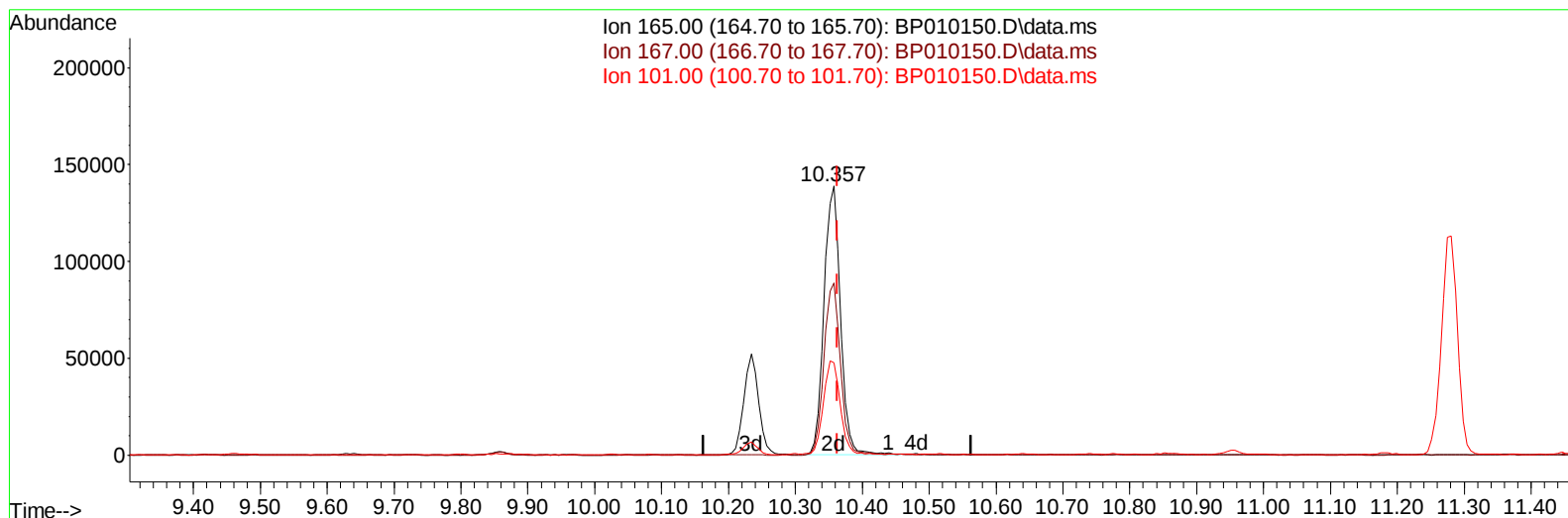
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 Sample : N2587-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP010150.D\data.ms

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

10.357min (-0.006) 25.22 ng/ul m

response 238640

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.98
101.00	49.80	34.31#
0.00	0.00	0.00

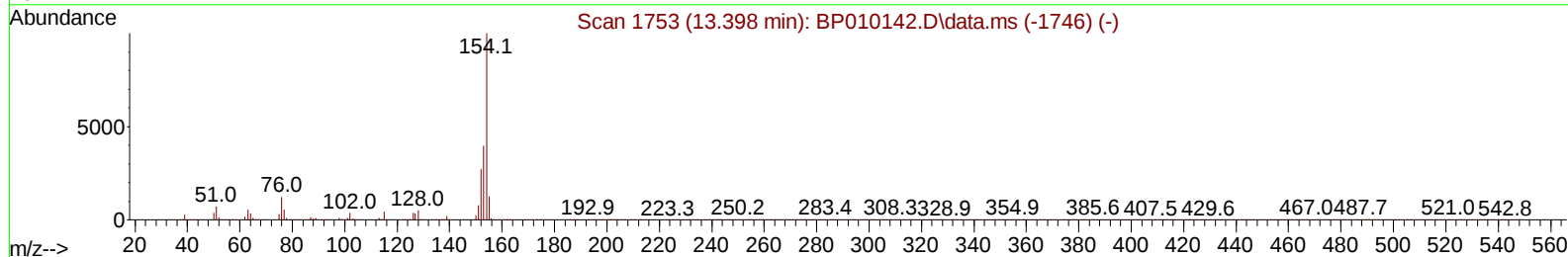
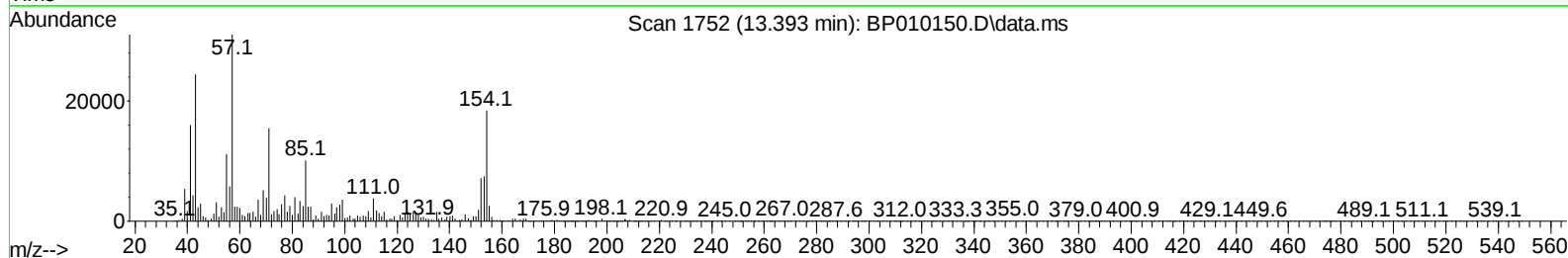
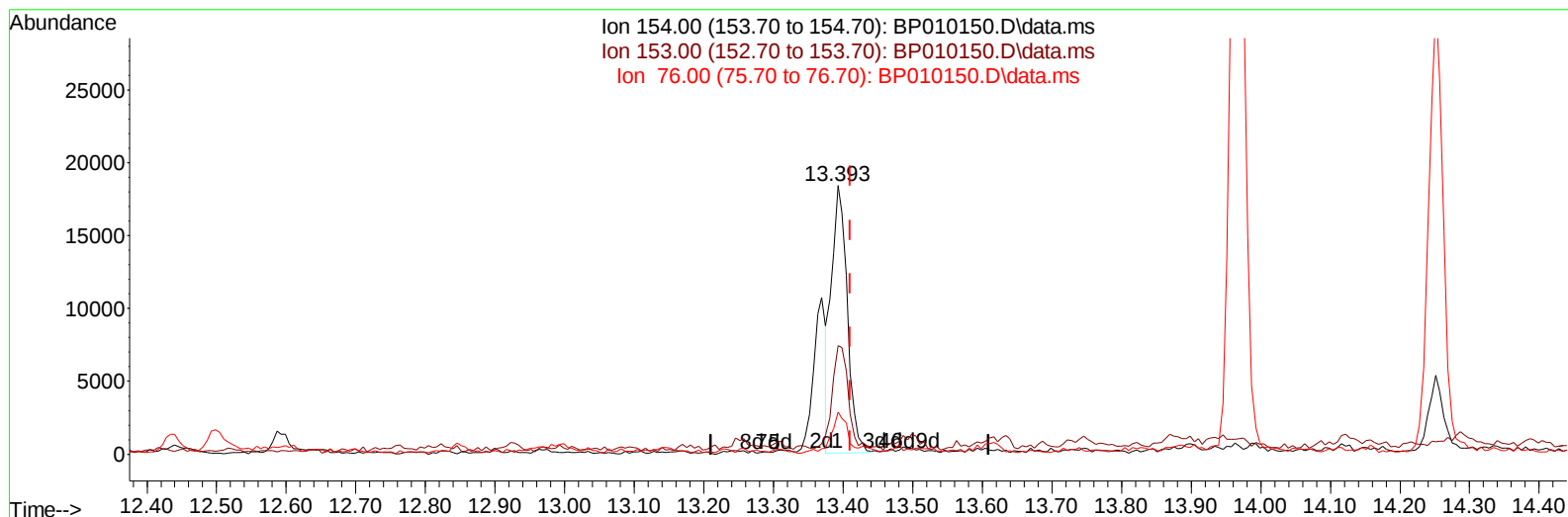
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Instrument :
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TIC: BP010150.D\data.ms

(43) 1,1'-Biphenyl

13.393min (-0.017) 1.15 ng/u1

response 28687

Ion	Exp%	Act%
154.00	100.00	100.00
153.00	35.80	40.41
76.00	10.60	15.65#
0.00	0.00	0.00

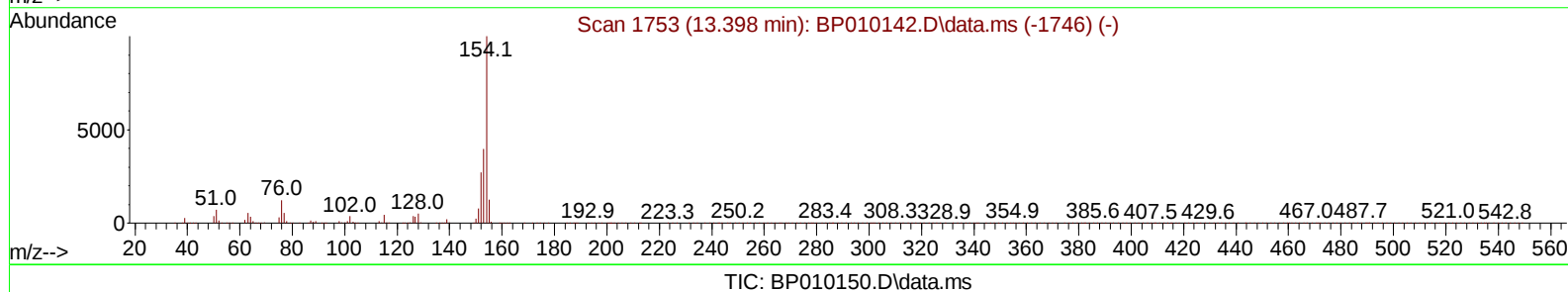
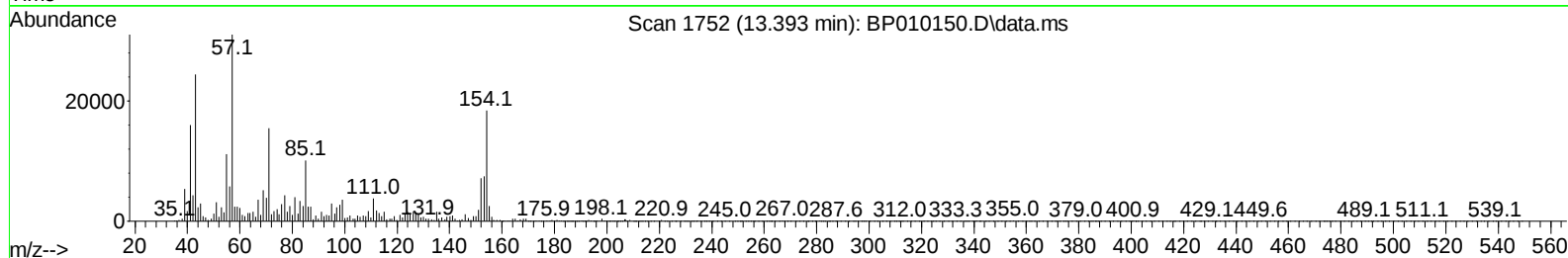
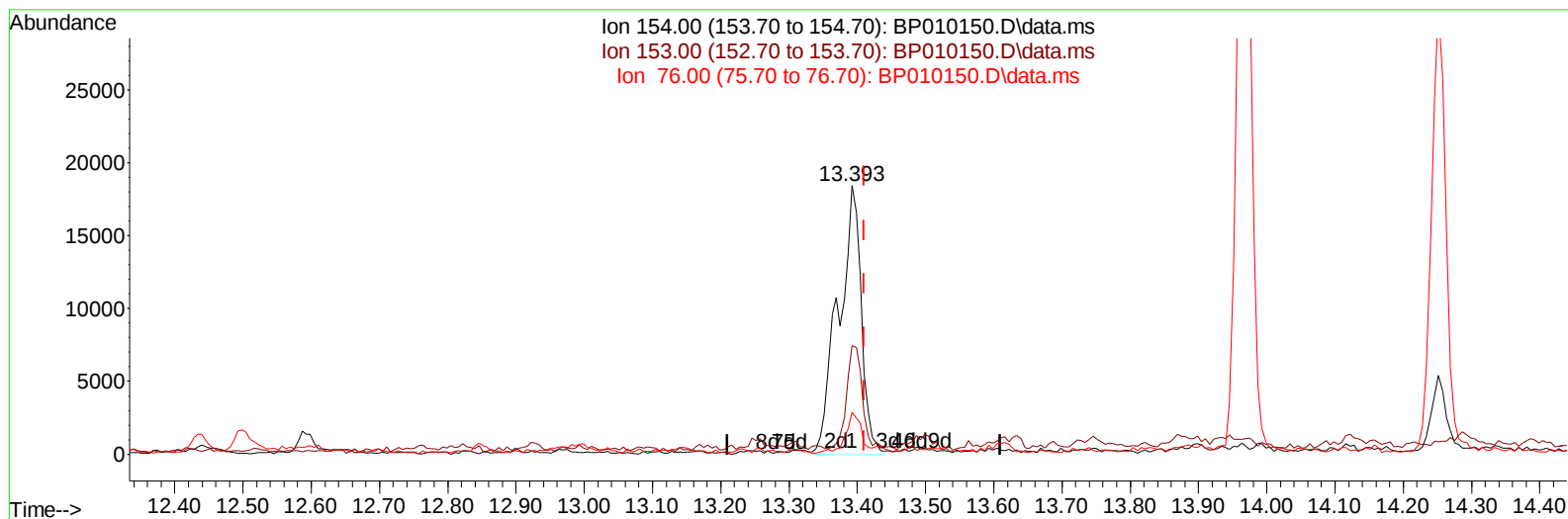
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 Acq On : 29 Apr 2022 14:28
 Operator : CG/JU
 Sample : N2587-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
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Reviewed By :Jagrut Upadhyay 05/02/2022
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(43) 1,1'-Biphenyl

13.393min (-0.017) 1.72 ng/ul m

response 42895

Ion	Exp%	Act%
154.00	100.00	100.00
153.00	35.80	40.41
76.00	10.60	15.65#
0.00	0.00	0.00

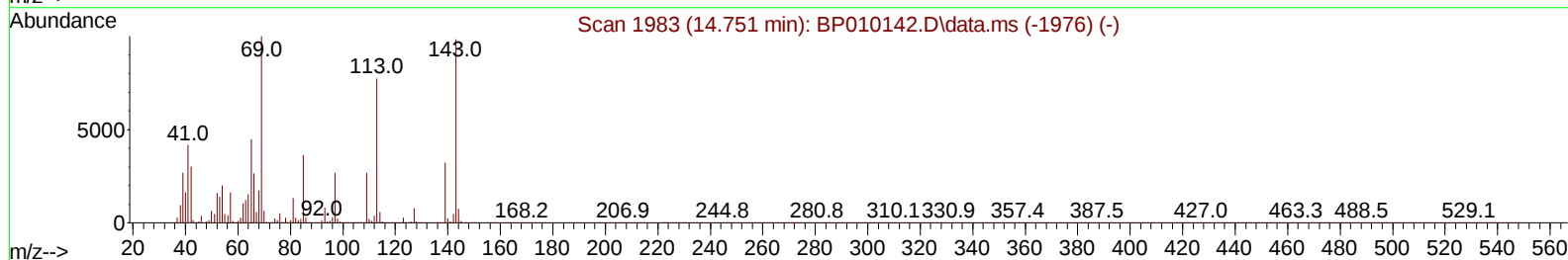
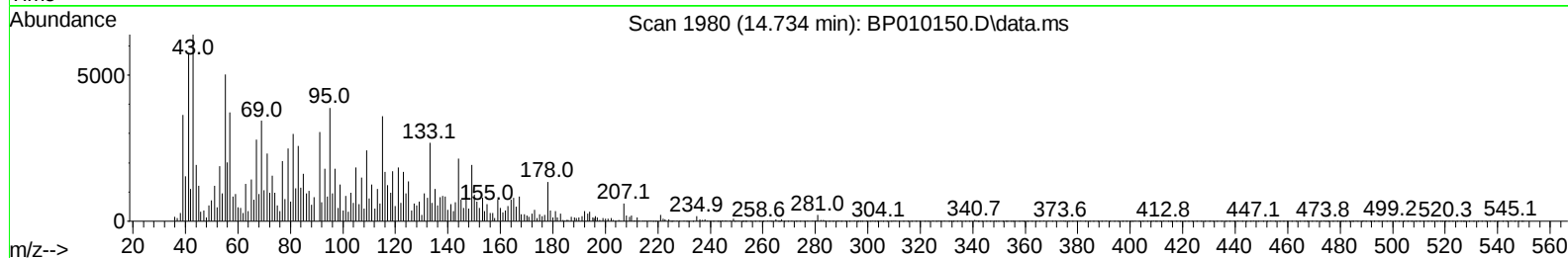
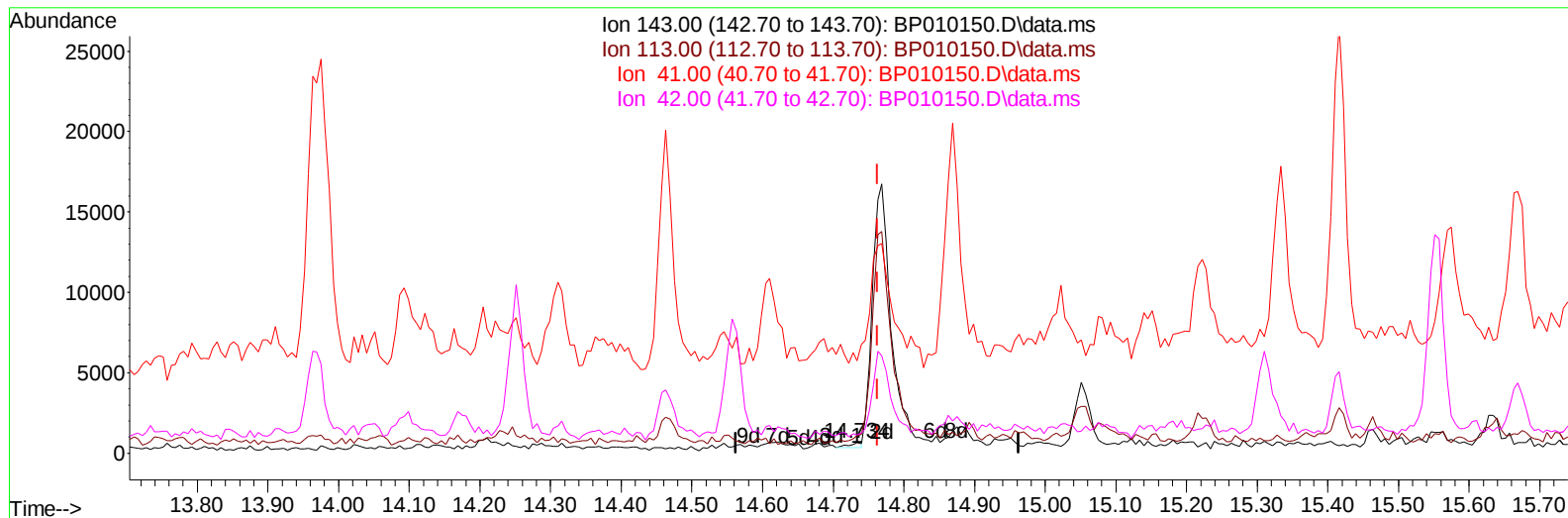
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TIC: BP010150.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.029) 0.08 ng/u1

response 371

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	169.66#
41.00	23.20	891.02#
42.00	18.00	172.45#

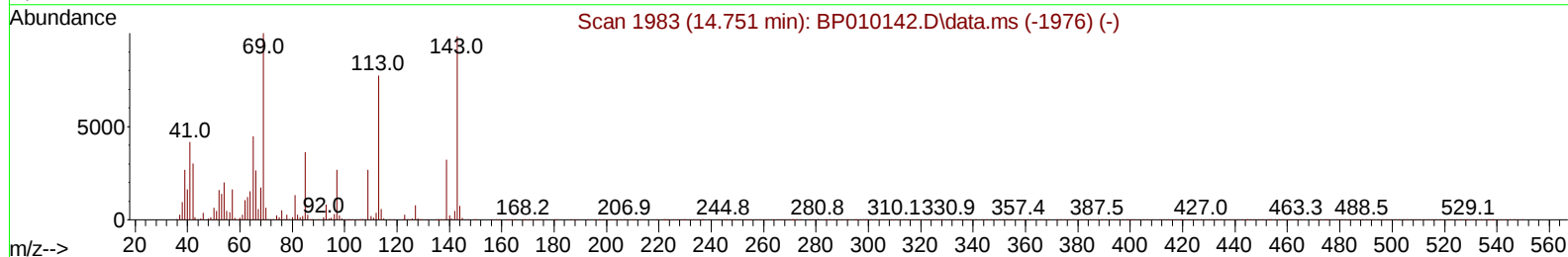
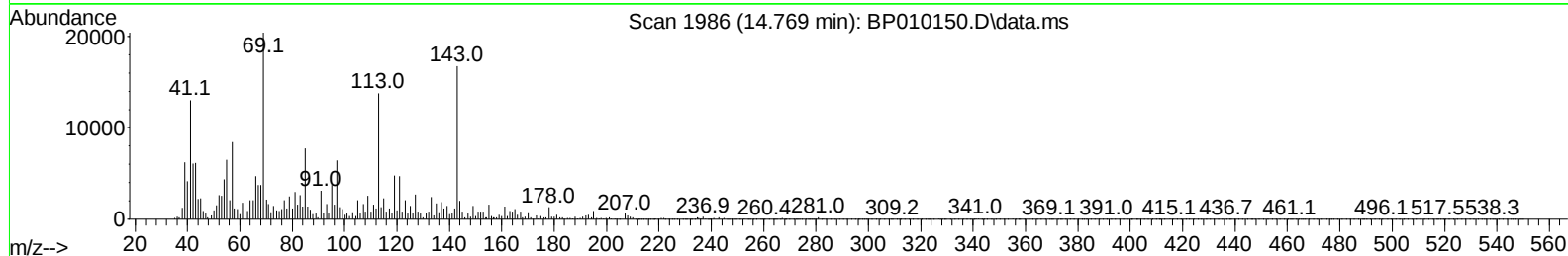
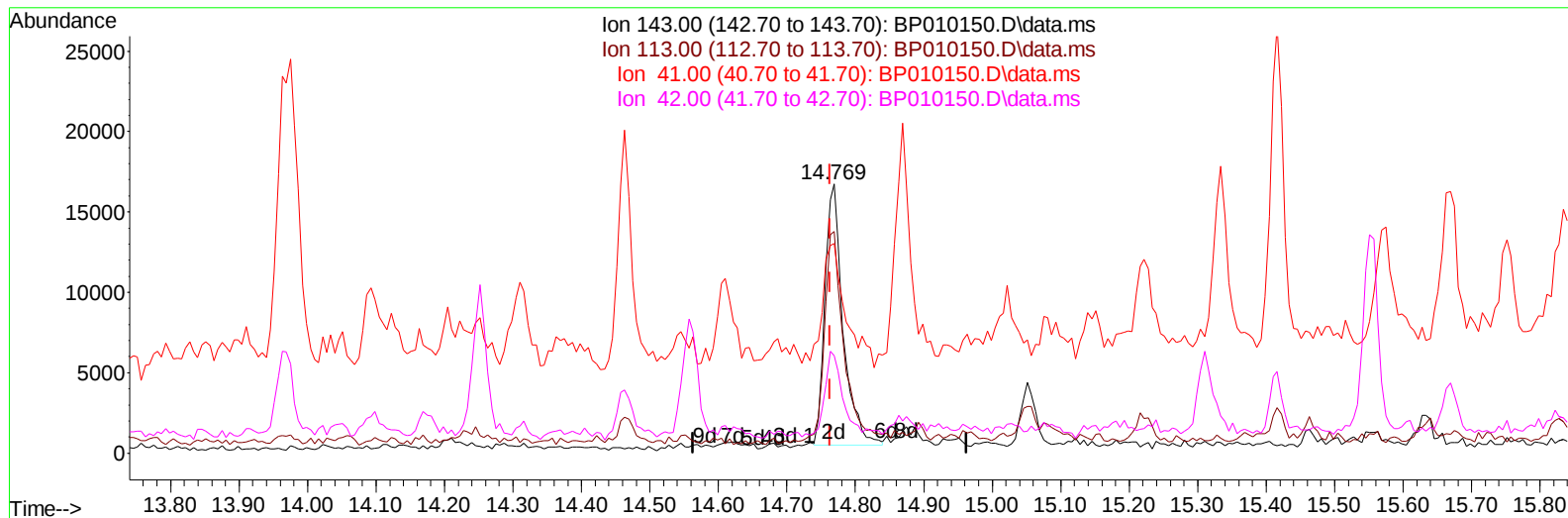
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Operator : CG/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
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EXET7

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TIC: BP010150.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 6.17 ng/ul m

response 29871

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.39#
41.00	23.20	77.88#
42.00	18.00	36.25#

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 Sample : N2587-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.928	152		145434	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.734	136		559659	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.557	164		337890	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.310	188		686504	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.398	240		670850	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264		710922	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.364	96		15823m	4.744	ng/uL	0.00
4) Pyridine-d5	3.787	84		75298	8.025	ng/ul	0.00
7) Phenol-d5	7.093	99		72419	5.474	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67		208516	26.269	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.458	132		210069	21.442	ng/ul	-0.01
15) 4-Methylphenol-d8	8.634	113		143567	13.099	ng/ul	-0.02
21) Nitrobenzene-d5	9.087	128		129619	29.120	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.810	143		142038	29.218	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165		238640m	25.216	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131		209304	15.424	ng/ul	-0.02
46) Dimethylphthalate-d6	13.969	166		842701	32.061	ng/ul	-0.01
49) Acenaphthylene-d8	14.251	160		943104	29.647	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143		29871m	6.173	ng/ul	0.00
60) Fluorene-d10	15.551	176		707242	31.280	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200		137695	31.721	ng/ul	-0.01
73) Anthracene-d10	17.410	188		1095425	32.982	ng/ul	-0.01
81) Pyrene-d10	19.645	212		1313348	34.425	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264		1262421	33.872	ng/ul	0.00
Target Compounds							
43) 1,1'-Biphenyl	13.393	154		42895m	1.721	ng/ul	Qvalue

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

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