

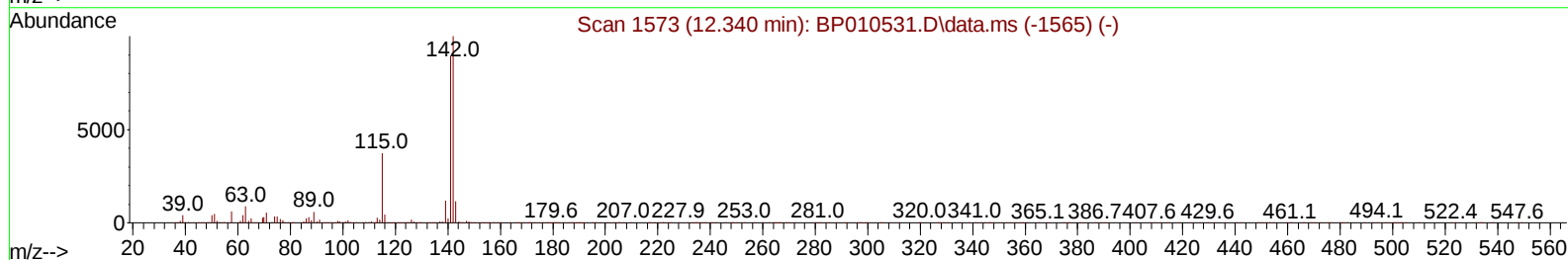
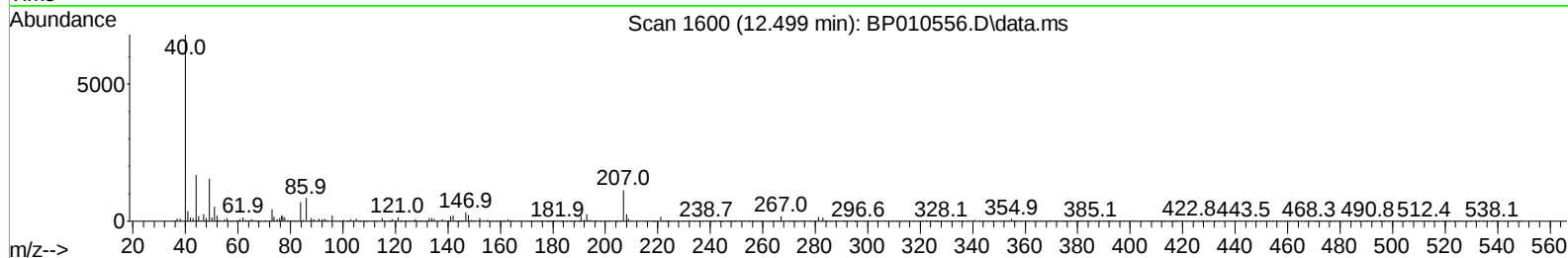
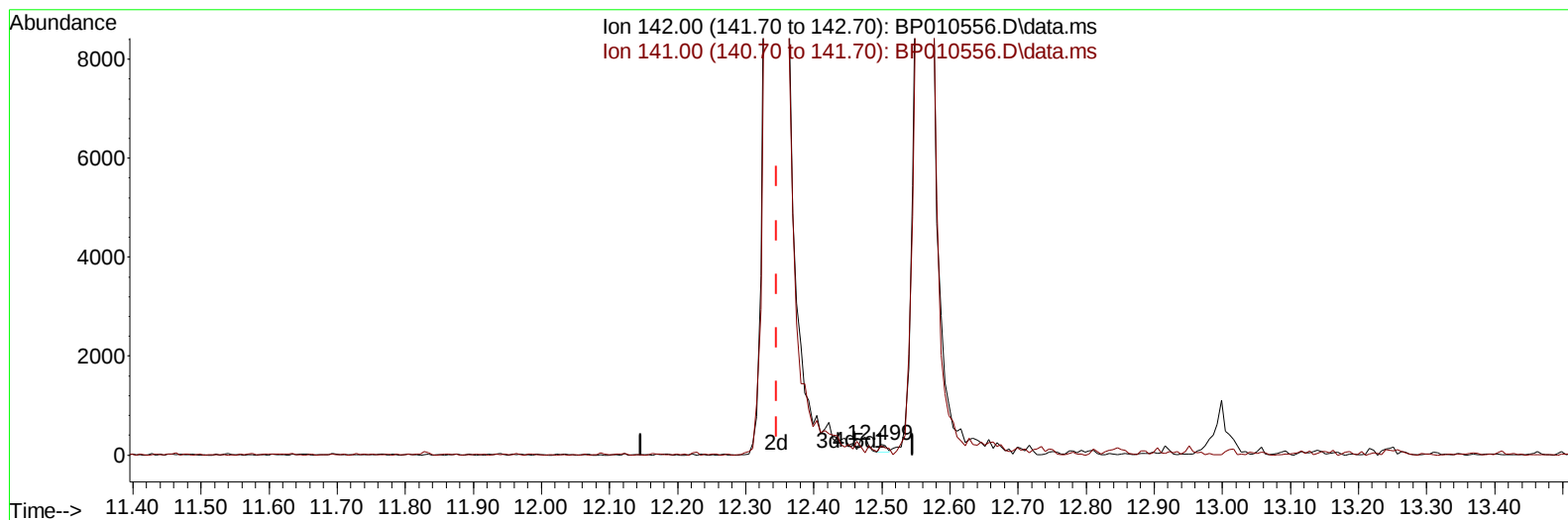
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010556.D
Acq On : 26 May 2022 02:35
Operator : CG/JU
Sample : N2950-09DL 10X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTH2DL

Manual IntegrationsAPPROVED

Quant Time: May 26 03:58:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 24 05:50:40 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/26/2022
Supervised By :mohammad ahmed 05/27/2022



TIC: BP010556.D\data.ms

(36) 2-Methylnaphthalene

12.499min (+ 0.153) 0.01 ng/ul

response 116

Ion	Exp%	Act%
142.00	100.00	100.00
141.00	80.40	80.26
0.00	0.00	0.00
0.00	0.00	0.00

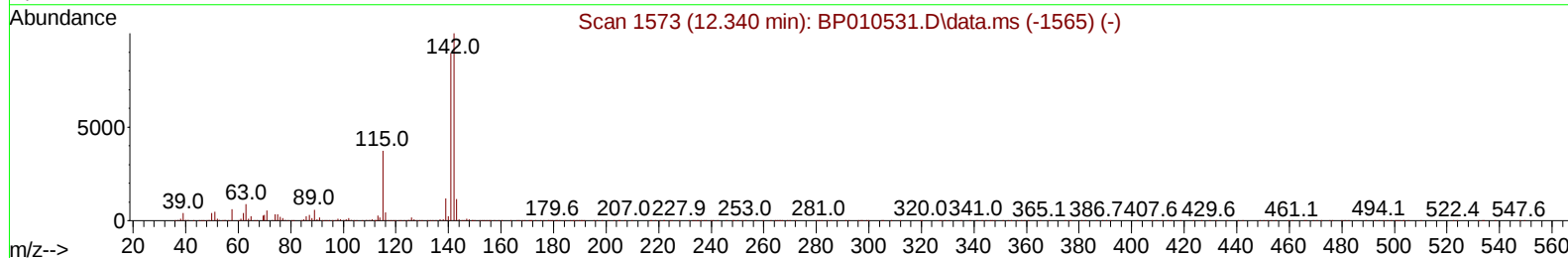
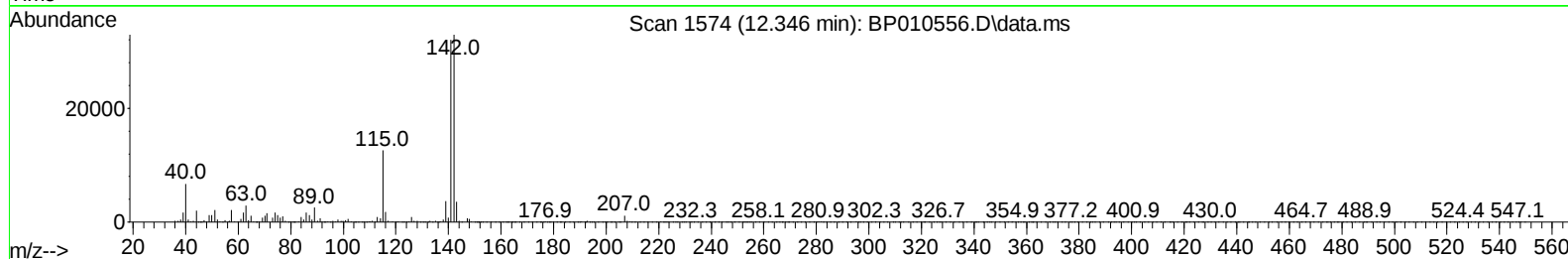
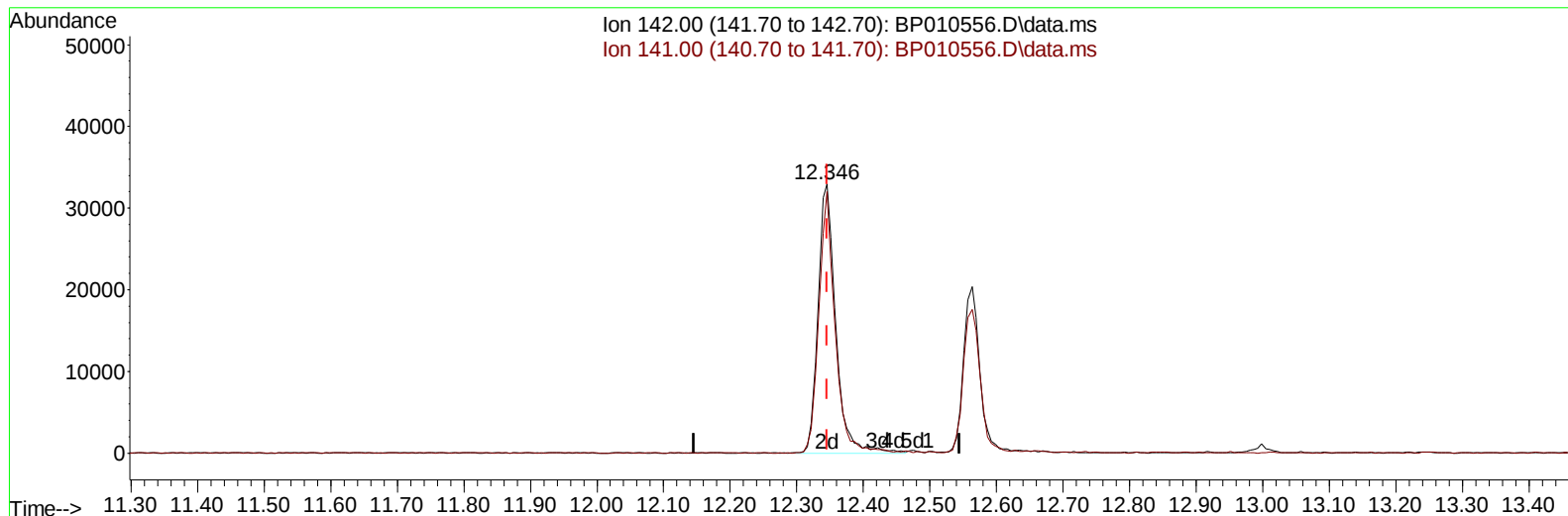
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TIC: BP010556.D\data.ms

(36) 2-Methylnaphthalene

12.346min (+ 0.000) 2.81 ng/ul m

response 60869

Ion	Exp%	Act%
142.00	100.00	100.00
141.00	80.40	97.28#
0.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	137469	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	585627	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164	417203	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.263	188	941702	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.345	240	908876	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	783376	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.058	99	6185	0.536	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	22270	2.873	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	19998	2.265	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	11126	1.135	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	11997	2.647	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	11175	2.327	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	10.834	131	23168	1.716	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	108623	3.534	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	107500	3.032	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.510	176	93236	3.487	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.363	188	158536	3.718	ng/ul	0.00
81) Pyrene-d10	19.598	212	193640	4.012	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	134474	3.419	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.728	128	1113220	36.394	ng/ul	96
36) 2-Methylnaphthalene	12.346	142	60869m	2.811	ng/ul	
37) 1-Methylnaphthalene	12.563	142	33636	1.538	ng/ul#	99
52) Acenaphthene	14.575	153	46819	1.871	ng/ul#	90
56) Dibenzofuran	14.916	168	84824	2.313	ng/ul	98
61) Fluorene	15.563	166	43074	1.436	ng/ul#	96
71) Pentachlorophenol	16.928	266	7282	1.040	ng/ul#	73
72) Phenanthrene	17.304	178	59734	1.189	ng/ul	99
77) Carbazole	17.669	167	170060	3.838	ng/ul#	95

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

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