

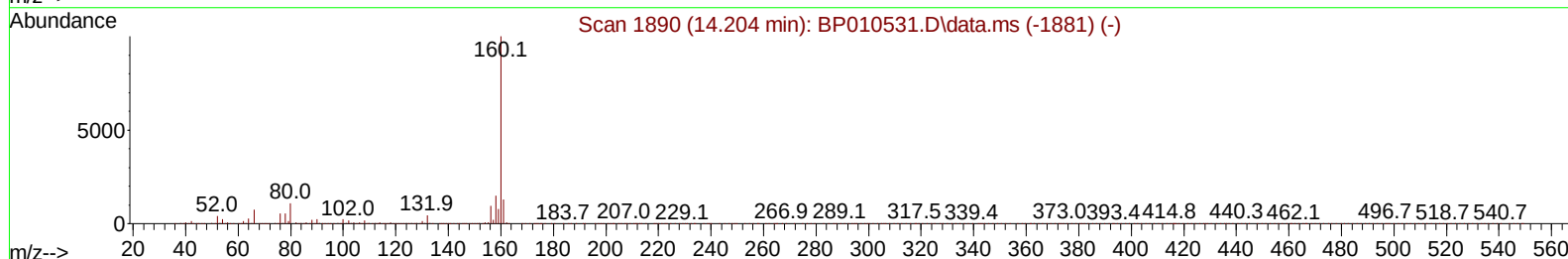
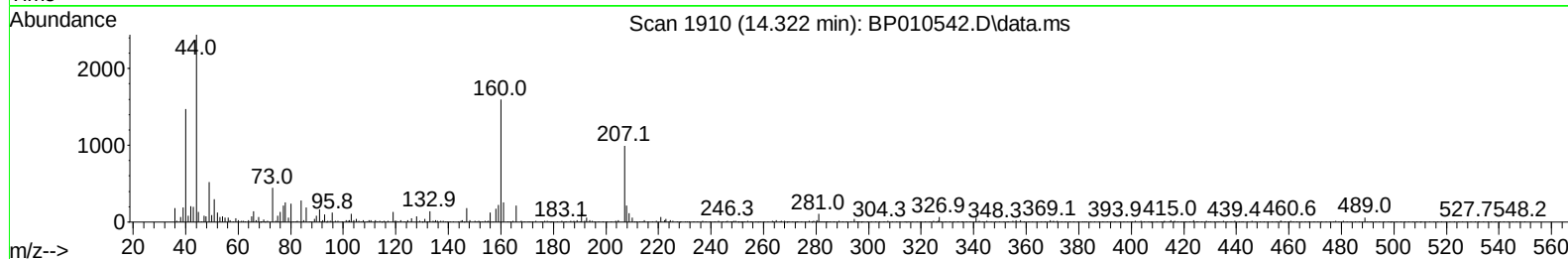
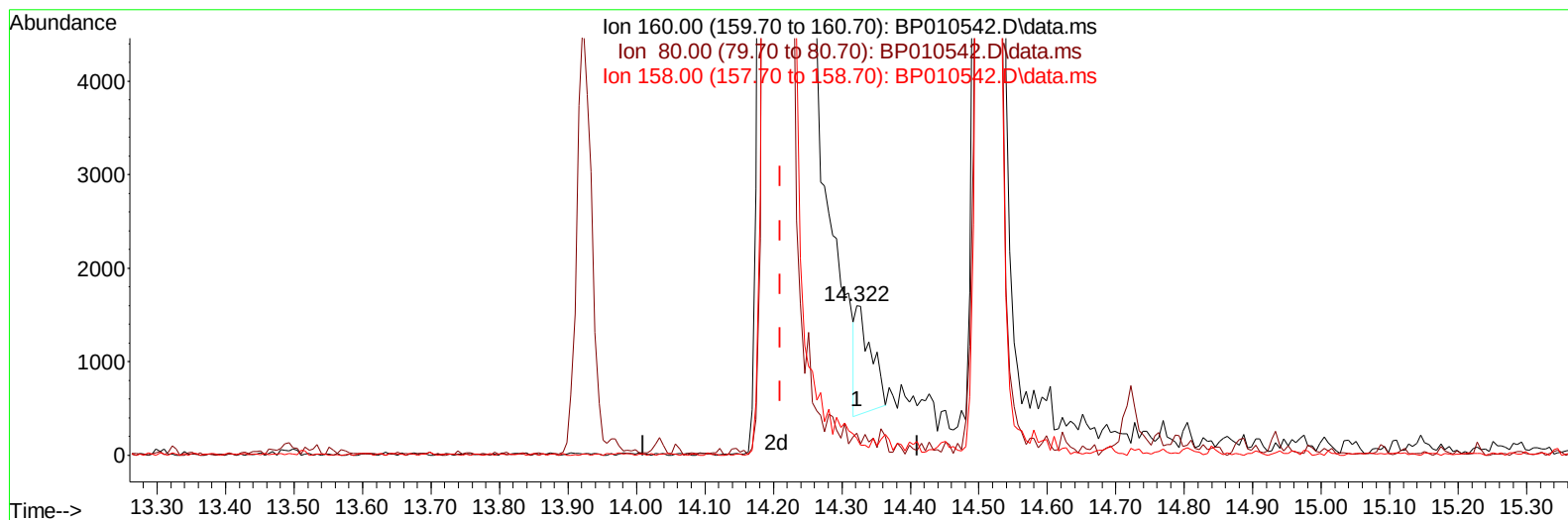
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
 Data File : BP010542.D
 Acq On : 25 May 2022 18:42
 Operator : CG/JU
 Sample : PB145042BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK042

Manual IntegrationsAPPROVED

Quant Time: May 25 23:27:41 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: BP010542.D\data.ms

(49) Acenaphthylene-d8 (S)

14.322min (+ 0.112) 0.06 ng/u1

response	1831	
Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	14.86
158.00	13.10	11.10
0.00	0.00	0.00

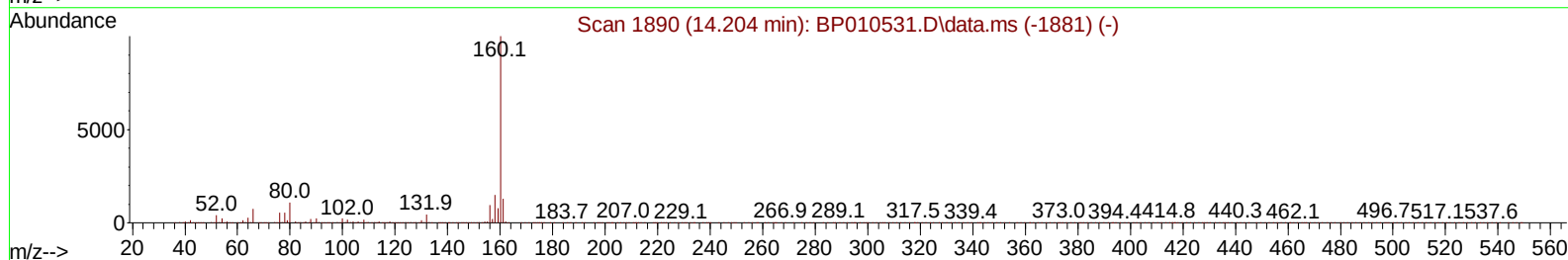
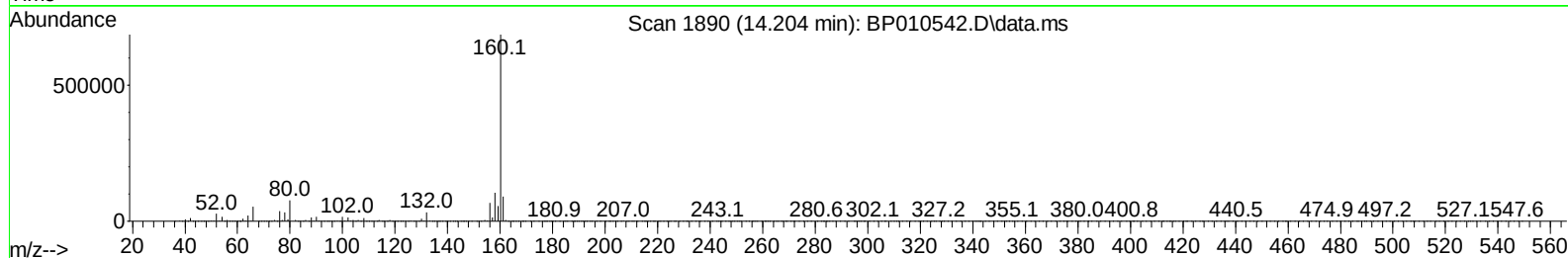
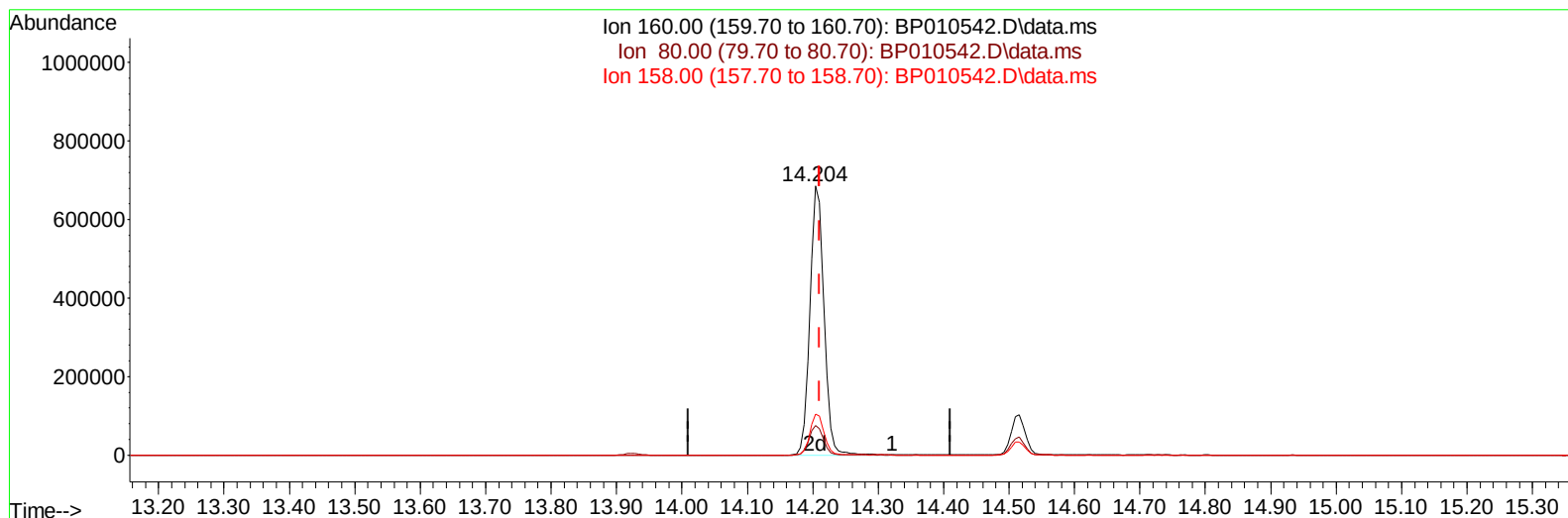
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TIC: BP010542.D\data.ms

(49) Acenaphthylene-d8 (S)

14.204min (-0.006) 35.44 ng/ul m

response 1040384

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	11.10#
158.00	13.10	15.13
0.00	0.00	0.00

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 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	111451	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	489112	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	345432	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	776856	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.351	240	738574	20.000	ng/ul	# 0.00
88) Perylene-d12	23.768	264	641944	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	17653	6.751	ng/uL	0.00
4) Pyridine-d5	3.752	84	250431	33.732	ng/ul	0.00
7) Phenol-d5	7.052	99	303144	32.407	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	221412	35.229	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	249330	34.833	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	256884	32.314	ng/ul	0.00
21) Nitrobenzene-d5	9.040	128	135055	35.678	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	142235	35.456	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	237311	31.285	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	356351	31.596	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	954384	37.500	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	1040384m	35.440	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	121203	26.756	ng/ul	0.00
60) Fluorene-d10	15.504	176	790453	35.700	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	123571	27.198	ng/ul	0.00
73) Anthracene-d10	17.369	188	1289838	36.667	ng/ul	0.00
81) Pyrene-d10	19.598	212	1589589	40.527	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1192545	36.996	ng/ul	0.00

Target Compounds Qvalue

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

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