

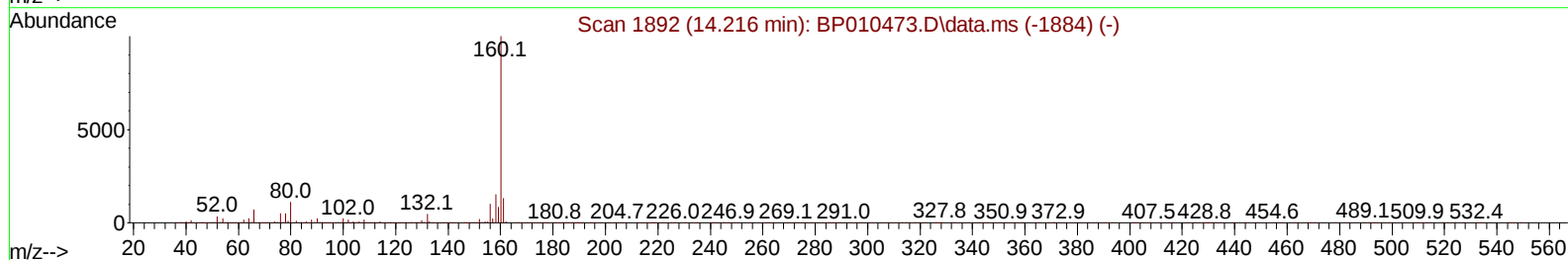
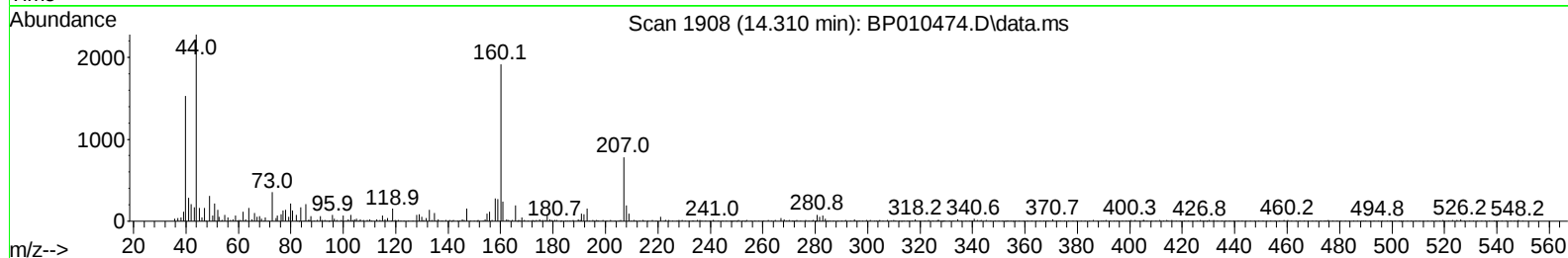
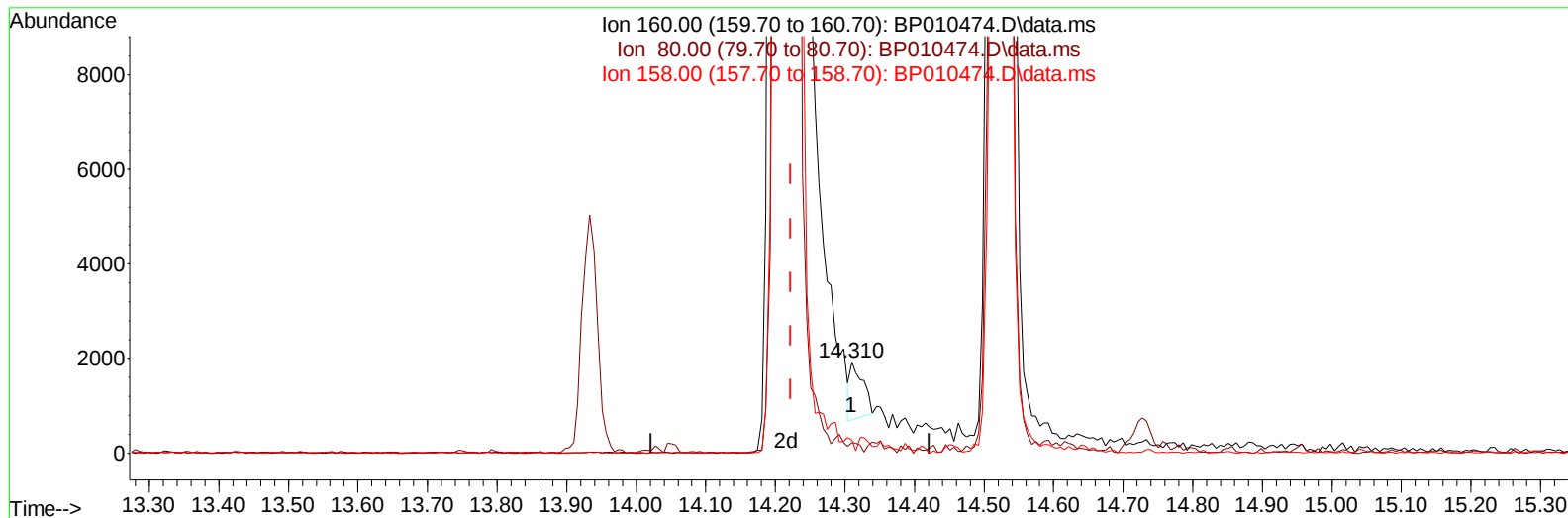
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
Data File : BP010474.D
Acq On : 23 May 2022 09:59
Operator : CG/JU
Sample : PB144961BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK961

Manual IntegrationsAPPROVED

Quant Time: May 23 23:18:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
Supervised By :Yogesh Patel 05/26/2022



TIC: BP010474.D\data.ms

(49) Acenaphthylene-d8 (S)

14.310min (+ 0.088) 0.04 ng/u1

response 1515

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	18.24
158.00	13.10	14.64
0.00	0.00	0.00

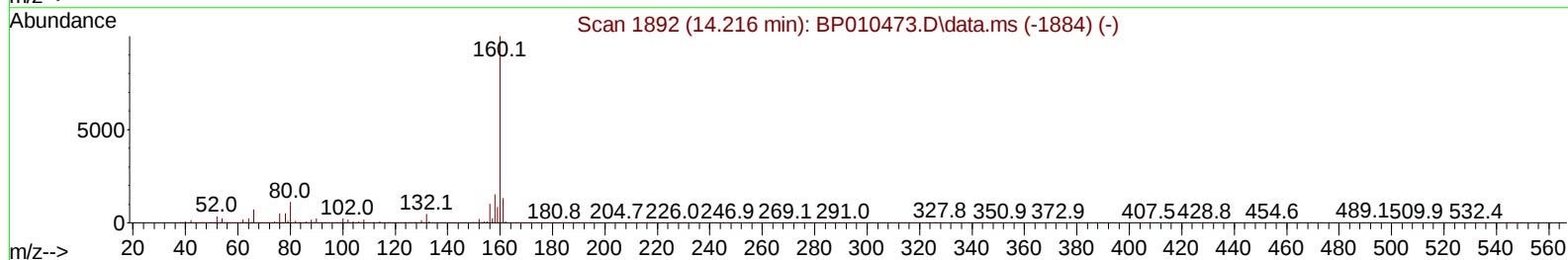
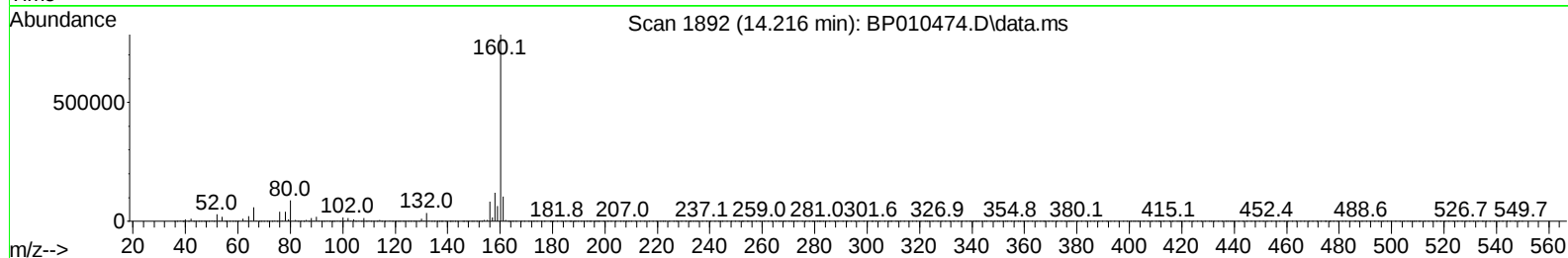
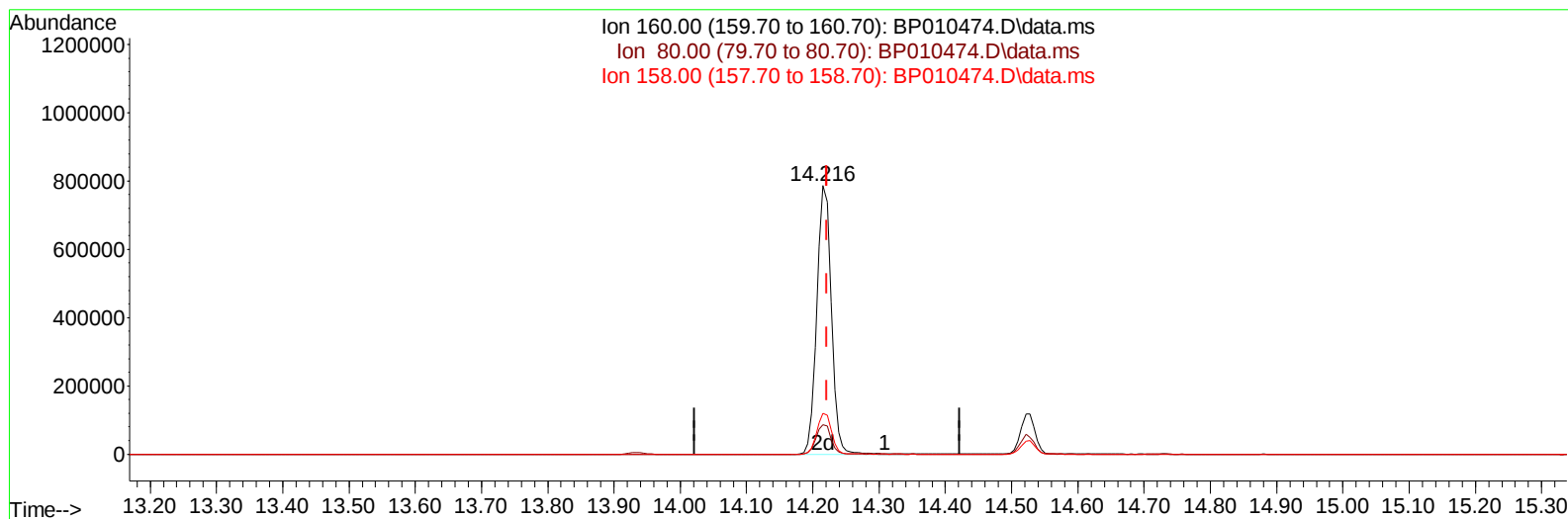
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TIC: BP010474.D\data.ms

(49) Acenaphthylene-d8 (S)

14.216min (-0.006) 33.99 ng/ul m

response 1190686

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	11.06#
158.00	13.10	15.17
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.892	152	135959	20.000	ng/ul	0.00
20) Naphthalene-d8	10.692	136	613145	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.521	164	412207	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.274	188	913390	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.356	240	865871	20.000	ng/ul	# 0.00
88) Perylene-d12	23.786	264	753618	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	18391	5.766	ng/uL	0.00
4) Pyridine-d5	3.757	84	267669	29.555	ng/ul	0.00
7) Phenol-d5	7.057	99	363394	31.845	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	245107	31.969	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	292686	33.519	ng/ul	-0.01
15) 4-Methylphenol-d8	8.604	113	315968	32.581	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	157904	33.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	170843	33.972	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	295300	31.055	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	431091	30.490	ng/ul	-0.01
46) Dimethylphthalate-d6	13.933	166	1022997	33.684	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1190686m	33.989	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	134141	24.815	ng/ul	0.00
60) Fluorene-d10	15.516	176	880988	33.344	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	130884	24.502	ng/ul	0.00
73) Anthracene-d10	17.374	188	1424775	34.448	ng/ul	0.00
81) Pyrene-d10	19.609	212	1753274	38.128	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1372920	36.280	ng/ul	0.00

Target Compounds Qvalue

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

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