

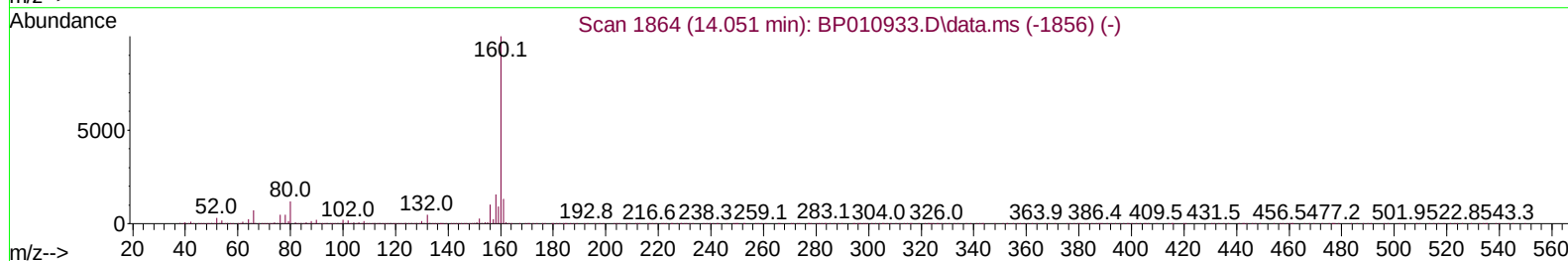
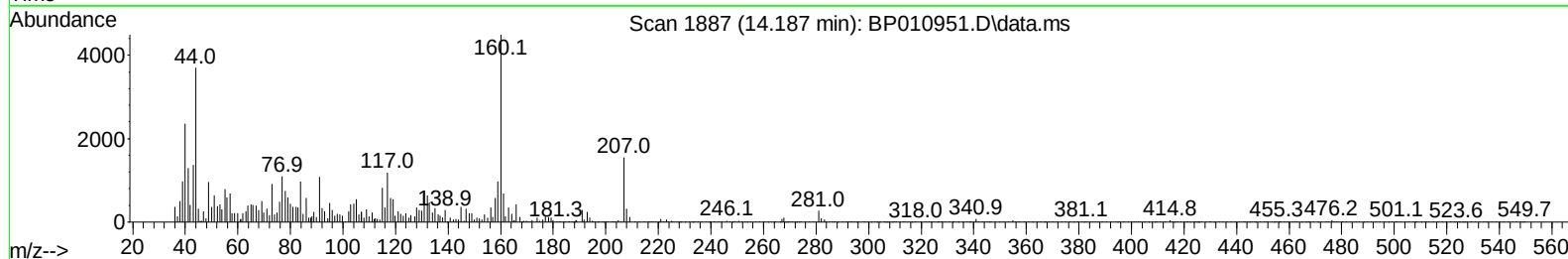
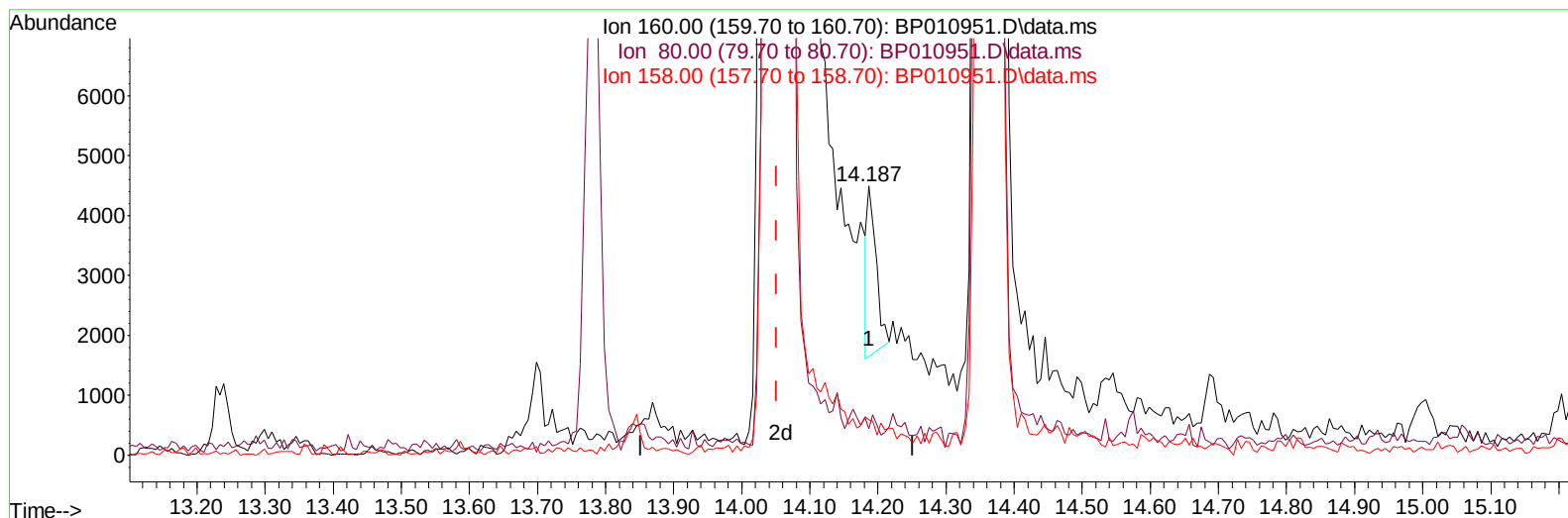
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071122\
Data File : BP010951.D
Acq On : 11 Jul 2022 16:25
Operator : CG/JU
Sample : N3623-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COAN5

Manual IntegrationsAPPROVED

Quant Time: Jul 11 18:16:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 06 23:13:53 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/12/2022
Supervised By :mohammad ahmed 07/12/2022



TIC: BP010951.D\data.ms

(49) Acenaphthylene-d8 (S)

14.187min (+ 0.135) 0.03 ng/u1

response 2570

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	10.10	9.73
158.00	15.30	13.02
0.00	0.00	0.00

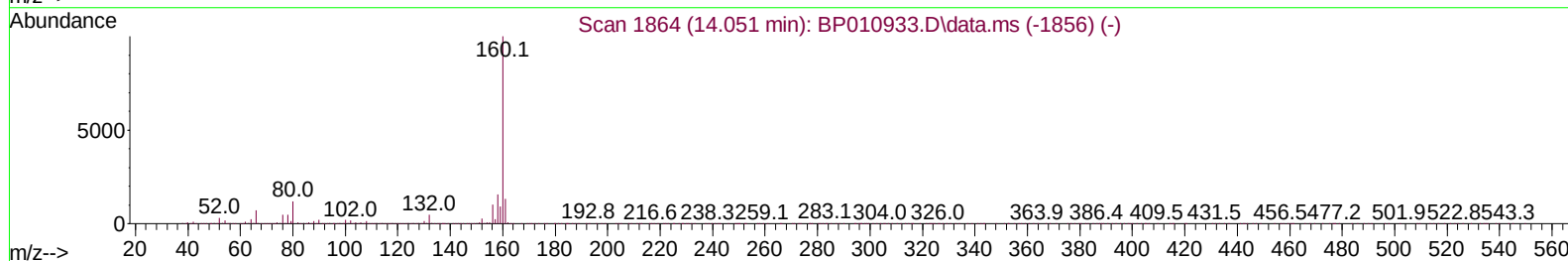
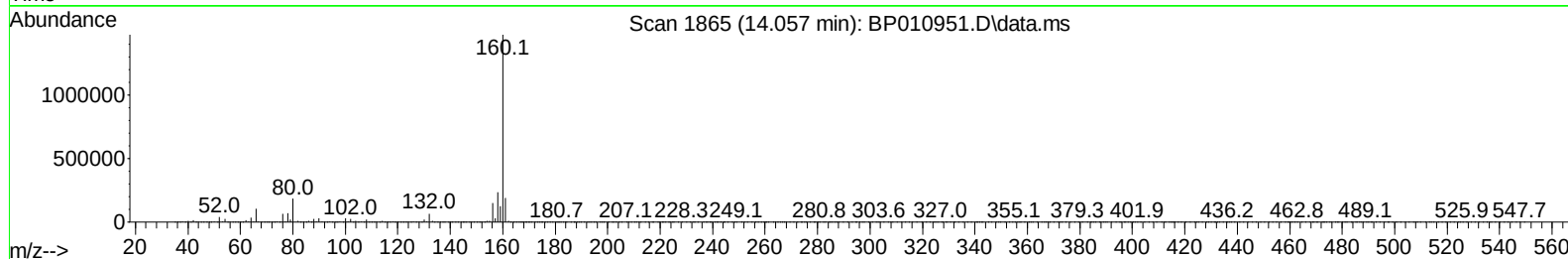
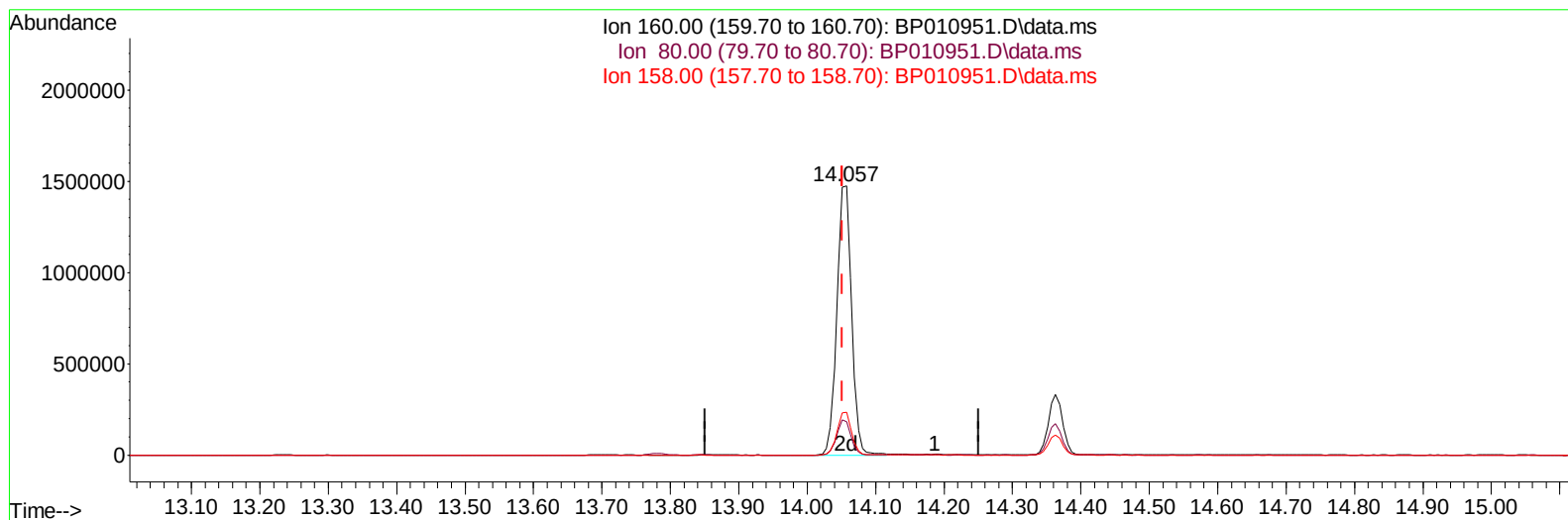
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(49) Acenaphthylene-d8 (S)

14.057min (+ 0.006) 24.17 ng/ul m

response 2209884

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	10.10	12.40#
158.00	15.30	15.88
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	7.705	152	514437	20.000	ng/ul	0.00
20) Naphthalene-d8	10.499	136	2053467	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	1045949	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.128	188	2046859	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.245	240	2096779	20.000	ng/ul	# 0.00
88) Perylene-d12	23.604	264	2291192	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	57496	4.228	ng/uL	0.00
4) Pyridine-d5	3.611	84	289674	8.002	ng/ul	0.00
7) Phenol-d5	6.881	99	186174	4.572	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	649012	24.369	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	621809	18.404	ng/ul	0.00
15) 4-Methylphenol-d8	8.416	113	353188	10.971	ng/ul	0.00
21) Nitrobenzene-d5	8.863	128	369637	28.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	322664	29.835	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	568558	21.276	ng/ul	0.00
31) 4-Chloroaniline-d4	10.646	131	472365	10.829	ng/ul	0.00
46) Dimethylphthalate-d6	13.781	166	1937718	27.045	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	2209884m	24.167	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	49547	4.427	ng/ul	0.00
60) Fluorene-d10	15.363	176	1725136	28.058	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	191648	28.261	ng/ul	0.00
73) Anthracene-d10	17.228	188	2814128	30.901	ng/ul	0.00
81) Pyrene-d10	19.480	212	3325253	29.109	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.457	264	3588983	31.699	ng/ul	0.01

Target Compounds Qvalue

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

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