

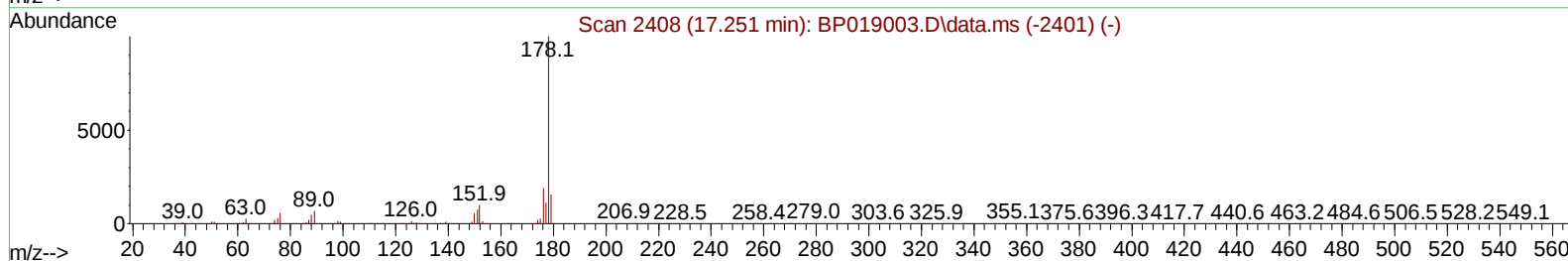
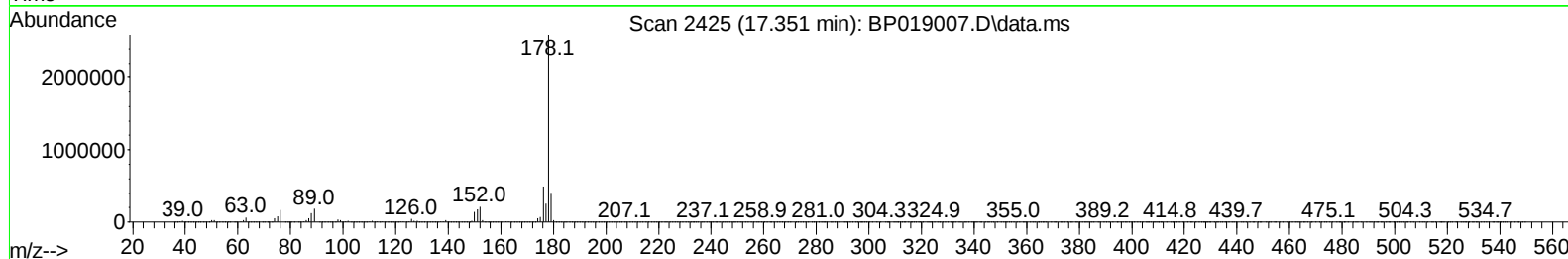
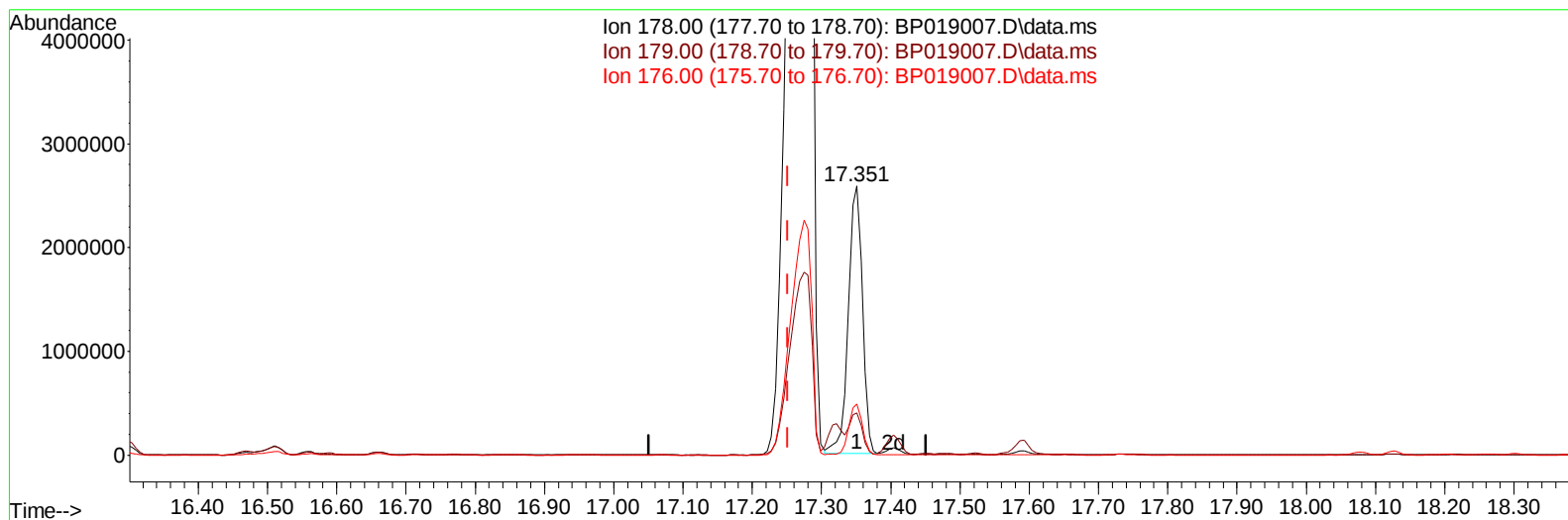
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019007.D
 Acq On : 21 Dec 2023 19:39
 Operator : MA/JU
 Sample : 05626-10ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCLF6ME

Manual IntegrationsAPPROVED

Quant Time: Dec 22 08:48:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/22/2023
 Supervised By :mohammad ahmed 12/25/2023



TIC: BP019007.D\data.ms

(72) Phenanthrene

17.351min (+ 0.100) 79.65 ng/u1

response 3620641

Ion	Exp%	Act%
178.00	100.00	100.00
179.00	15.40	15.61
176.00	19.30	18.98
0.00	0.00	0.00

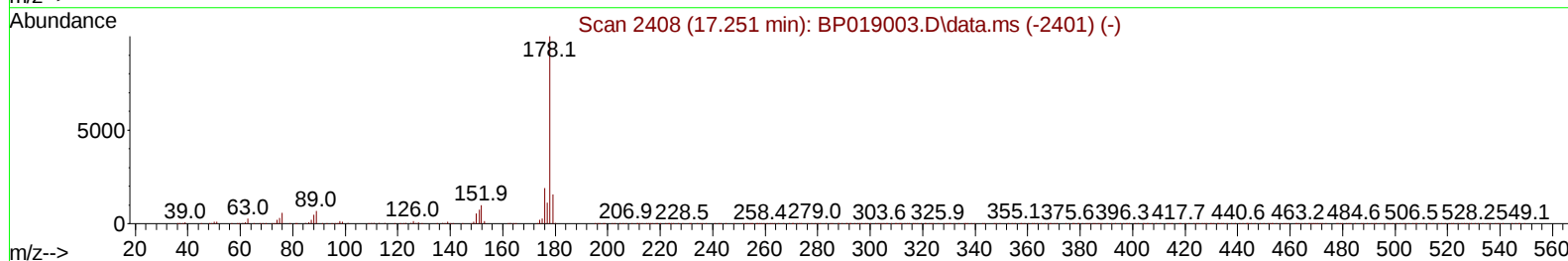
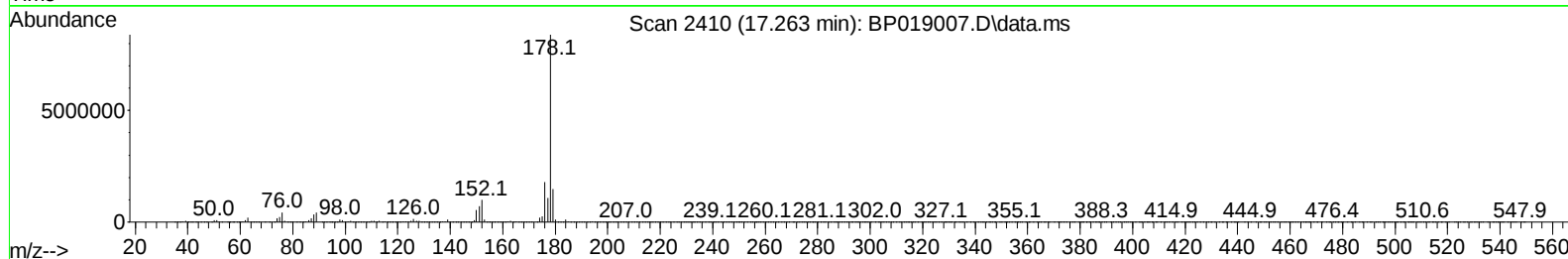
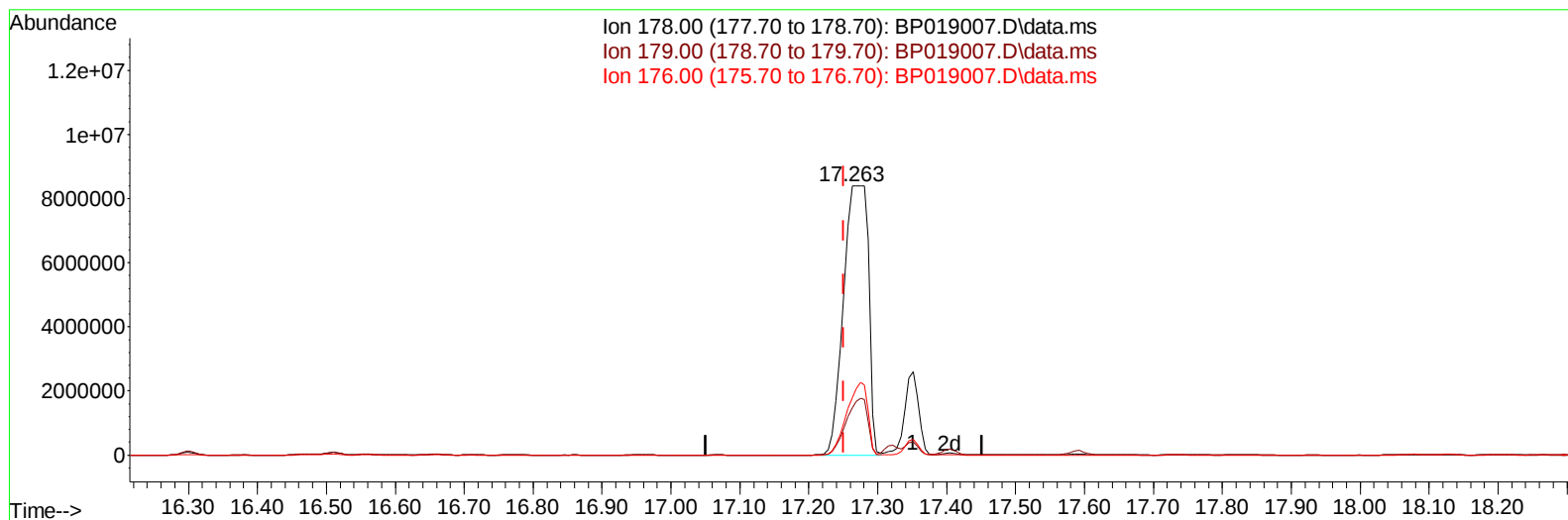
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019007.D
 Acq On : 21 Dec 2023 19:39
 Operator : MA/JU
 Sample : 05626-10ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCLF6ME

Manual IntegrationsAPPROVED

Quant Time: Dec 22 08:48:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/22/2023
 Supervised By :mohammad ahmed 12/25/2023



TIC: BP019007.D\data.ms

(72) Phenanthrene

17.263min (+ 0.012) 464.97 ng/u1 m

response 21135802

Ion	Exp%	Act%
178.00	100.00	100.00
179.00	15.40	17.52
176.00	19.30	21.35
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019007.D
 Acq On : 21 Dec 2023 19:39
 Operator : MA/JU
 Sample : 05626-10ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCLF6ME

Manual IntegrationsAPPROVED

Quant Time: Dec 22 23:27:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/22/2023
 Supervised By :mohammad ahmed 12/25/2023

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.822	152		130360	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136		502753	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164		339712	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188		752645	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.304	240		801176	20.000	ng/ul	# 0.00
88) Perylene-d12	23.657	264		972490	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.252	96		3194	0.835	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	6.999	99		59288	4.532	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67		40272	5.194	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132		50547	4.952	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113		35419	3.345	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128		27445	6.133	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143		29377	6.332	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165		56791	5.987	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131		54129	4.366	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166		217908	7.187	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160		222489	6.664	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143		25968	5.218	ng/ul	0.01
60) Fluorene-d10	15.451	176		192667	7.570	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200		26799	5.938	ng/ul	0.02
73) Anthracene-d10	17.316	188		290181	7.334	ng/ul	0.01
81) Pyrene-d10	19.545	212		252099	4.042	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264		164806	2.906	ng/ul	0.00
Target Compounds							
						Qvalue	
30) Naphthalene	10.675	128		6218517	203.893	ng/ul	99
36) 2-Methylnaphthalene	12.281	142		3593457	166.312	ng/ul	100
37) 1-Methylnaphthalene	12.499	142		1843171	84.471	ng/ul	99
50) Acenaphthylene	14.175	152		139061	3.705	ng/ul#	88
52) Acenaphthene	14.534	153		6675208	268.960	ng/ul	99
61) Fluorene	15.516	166		6954326	254.790	ng/ul	99
72) Phenanthrene	17.263	178		21135802m	464.970	ng/ul	
74) Anthracene	17.351	178		3620641	79.463	ng/ul	100
80) Fluoranthene	19.233	202		14697867	199.379	ng/ul#	91
82) Pyrene	19.580	202		6884301	92.570	ng/ul#	90
85) Benzo(a)anthracene	21.286	228		3145436	48.369	ng/ul	99
87) Chrysene	21.339	228		2458789	41.058	ng/ul	97
90) Benzo(b)fluoranthene	22.939	252		1636811	22.828	ng/ul	97
91) Benzo(k)fluoranthene	22.986	252		521532	7.517	ng/ul	97
93) Benzo(a)pyrene	23.551	252		270079	4.122	ng/ul#	95
94) Indeno(1,2,3-cd)pyrene	26.098	276		138747	1.772	ng/ul#	93
95) Dibenzo(a,h)anthracene	26.121	278		79579	1.222	ng/ul	95

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019007.D
Acq On : 21 Dec 2023 19:39
Operator : MA/JU
Sample : 05626-10ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCLF6ME

Manual IntegrationsAPPROVED

Quant Time: Dec 22 23:27:03 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 21 04:46:12 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/22/2023
Supervised By :mohammad ahmed 12/25/2023

