

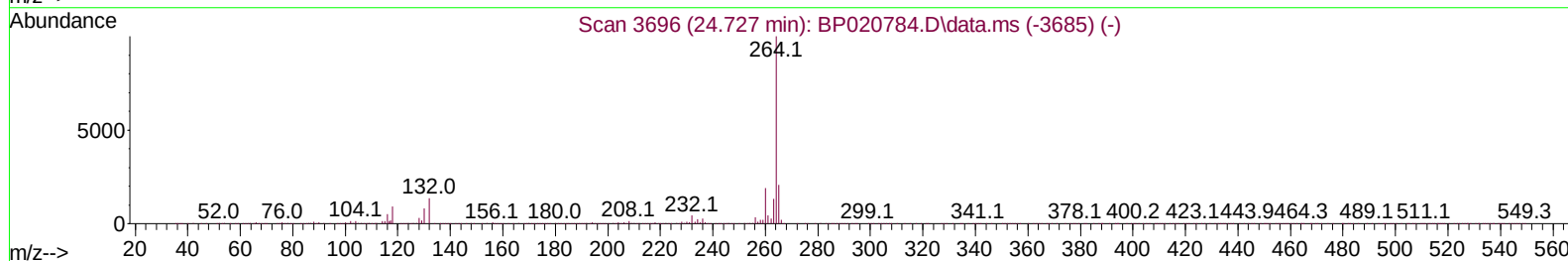
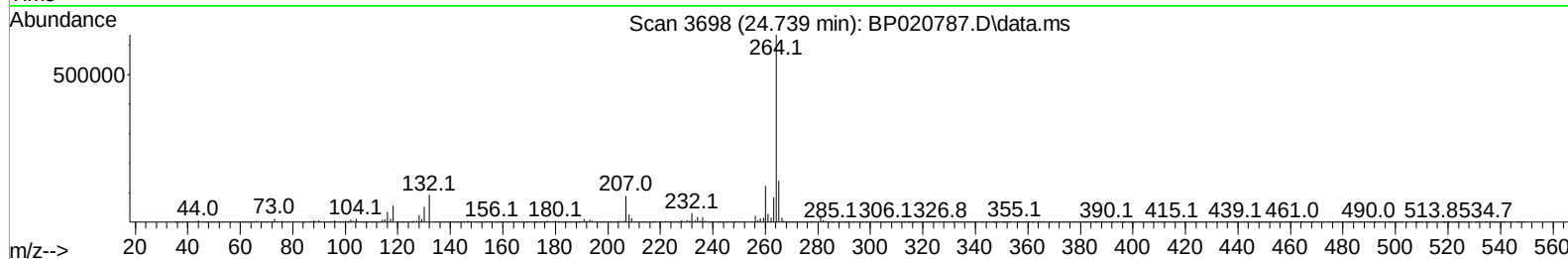
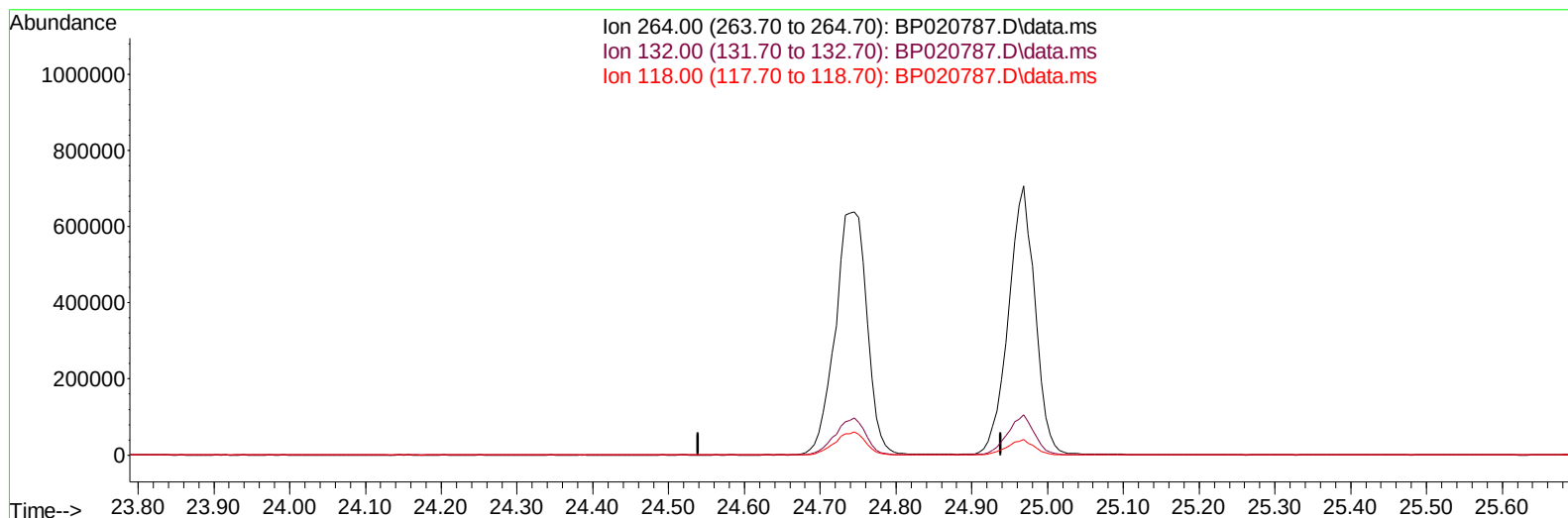
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020787.D
Acq On : 05 Jul 2024 13:40
Operator : MA/JU
Sample : P2987-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0T86

Manual IntegrationsAPPROVED

Quant Time: Jul 05 14:57:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 03 15:01:56 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/08/2024
Supervised By :mohammad ahmed 07/08/2024



TIC: BP020787.D\data.ms

(92) Benzo(a)pyrene-d12 (S)

24.739min (-24.739) 0.00 ng/u1

response 0

Ion	Exp%	Act%
264.00	100.00	0.00
132.00	15.50	0.00#
118.00	9.80	0.00#
0.00	0.00	0.00

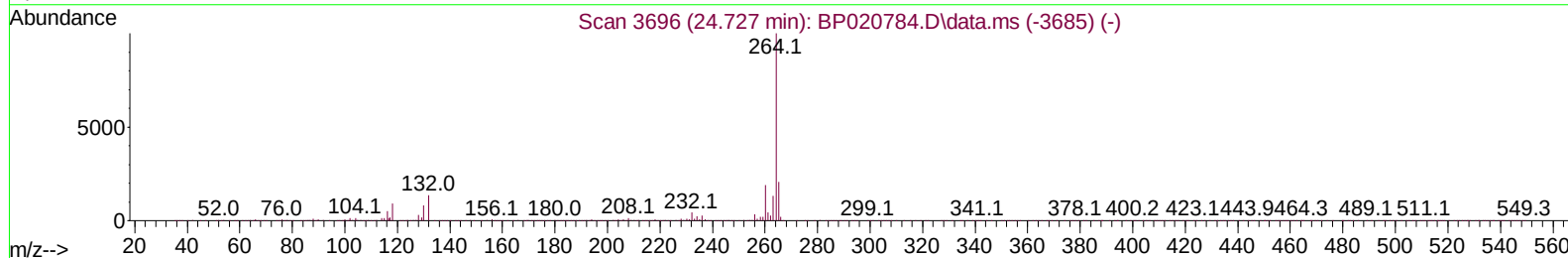
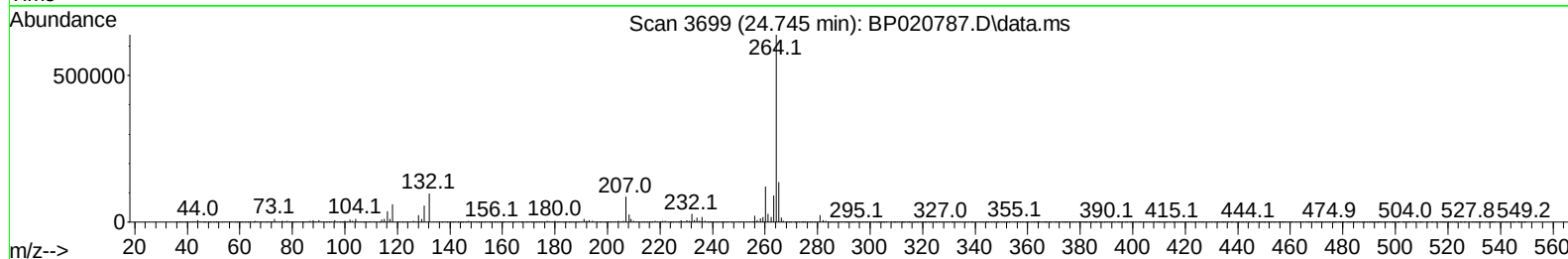
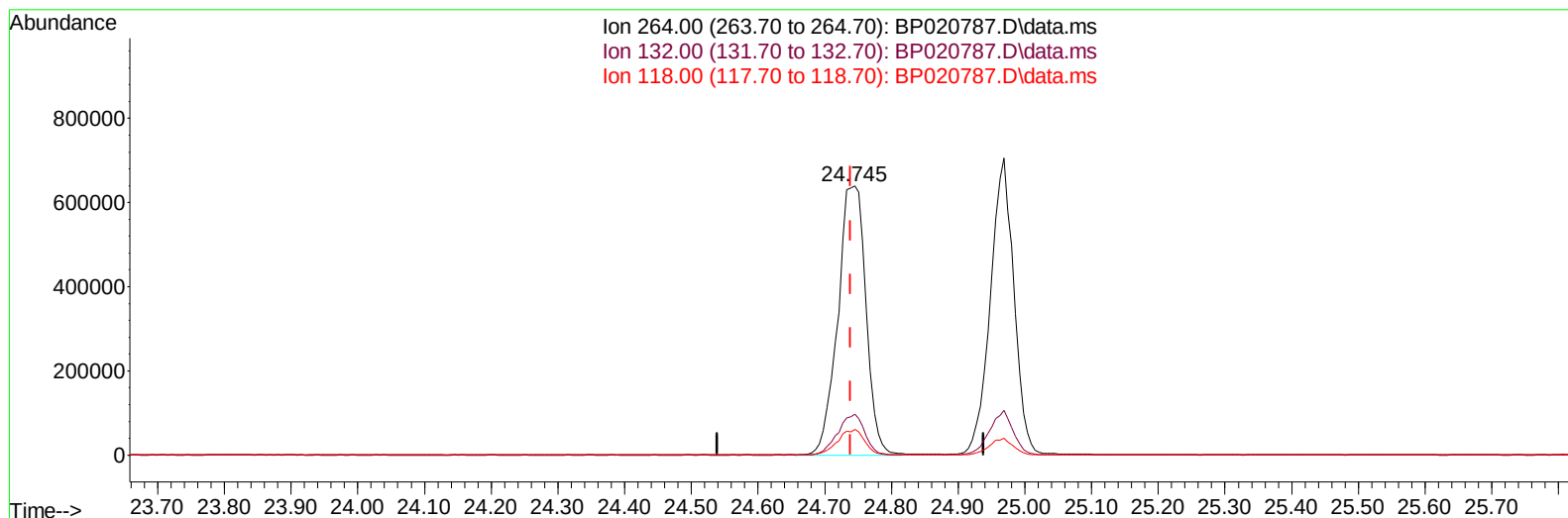
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(92) Benzo(a)pyrene-d12 (S)

24.745min (+ 0.006) 22.36 ng/ul m

response 1874217

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	15.50	15.26
118.00	9.80	9.67
0.00	0.00	0.00

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 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0T86

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/08/2024
 Supervised By :mohammad ahmed 07/08/2024

Quant Time: Jul 05 14:58:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 15:01:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	278172	20.000	ng/ul	0.00
20) Naphthalene-d8	10.493	136	1176096	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.357	164	770163	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.169	188	1642381	20.000	ng/ul	0.00
79) Chrysene-d12	21.616	240	1536851	20.000	ng/ul	0.00
88) Perylene-d12	24.968	264	1723436	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	20244	3.125	ng/uL	0.00
4) Pyridine-d5	3.675	84	220939	11.703	ng/ul	0.00
7) Phenol-d5	6.916	99	477714	20.938	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	277123	19.160	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.269	132	394601	21.162	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	397151	21.433	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	196202	22.596	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.587	143	222242	25.576	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	392097	21.906	ng/ul	0.00
31) 4-Chloroaniline-d4	10.640	131	363982	14.510	ng/ul	0.00
46) Dimethylphthalate-d6	13.769	166	1224186	22.861	ng/ul	0.01
49) Acenaphthylene-d8	14.045	160	1380772	21.609	ng/ul	0.00
54) 4-Nitrophenol-d4	14.592	143	179407	21.928	ng/ul	0.01
60) Fluorene-d10	15.369	176	1034308	22.954	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	149907	18.112	ng/ul	0.00
73) Anthracene-d10	17.275	188	1629263	22.104	ng/ul	0.01
81) Pyrene-d10	19.651	212	1945363	21.070	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.745	264	1874217m	22.356	ng/ul	0.00

Target Compounds Qvalue

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

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