

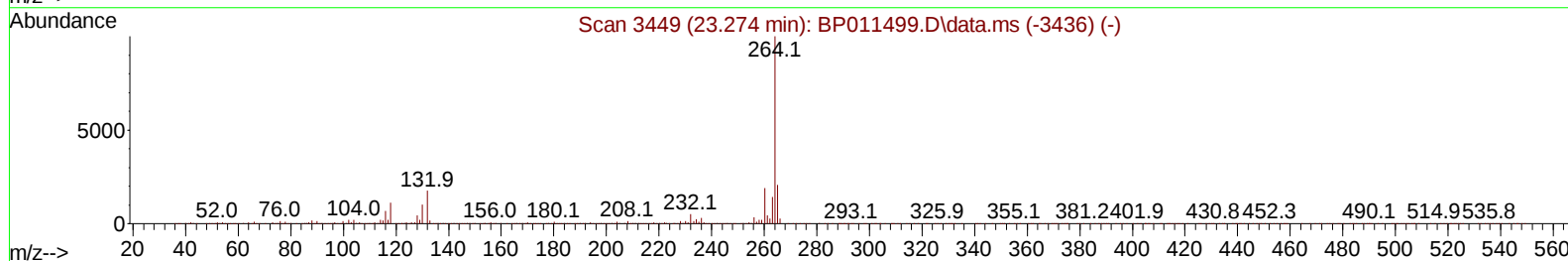
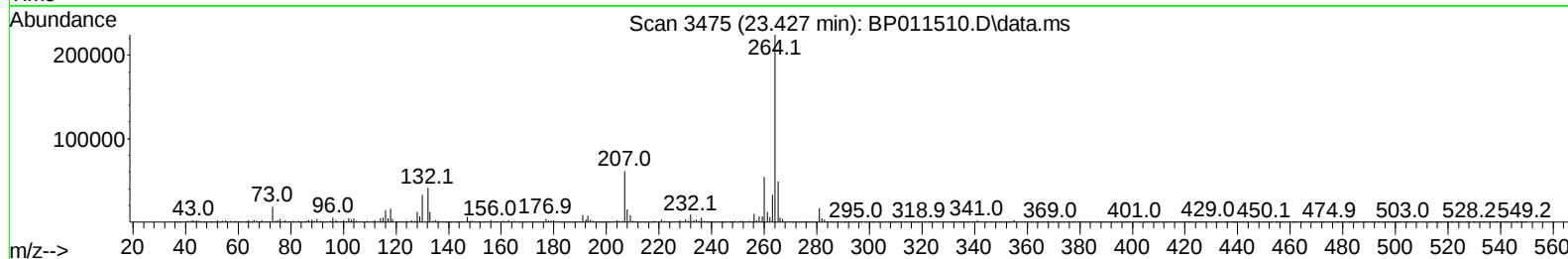
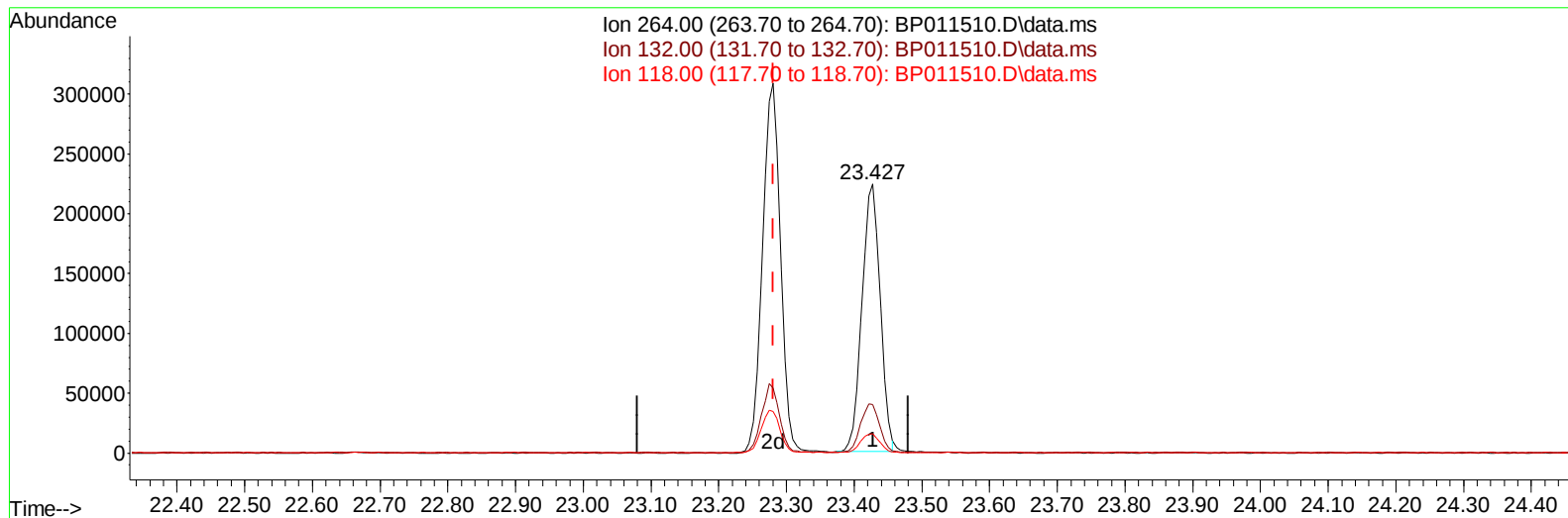
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011510.D
Acq On : 19 Aug 2022 16:37
Operator : CG/JU
Sample : N4221-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KQ0

Manual IntegrationsAPPROVED

Quant Time: Aug 20 04:38:08 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 01:58:37 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/22/2022
Supervised By :Sohil Jodhani 08/22/2022



TIC: BP011510.D\data.ms

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

23.427min (+ 0.147) 20.42 ng/u1

response 420218

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	17.60	18.17
118.00	11.10	8.99
0.00	0.00	0.00

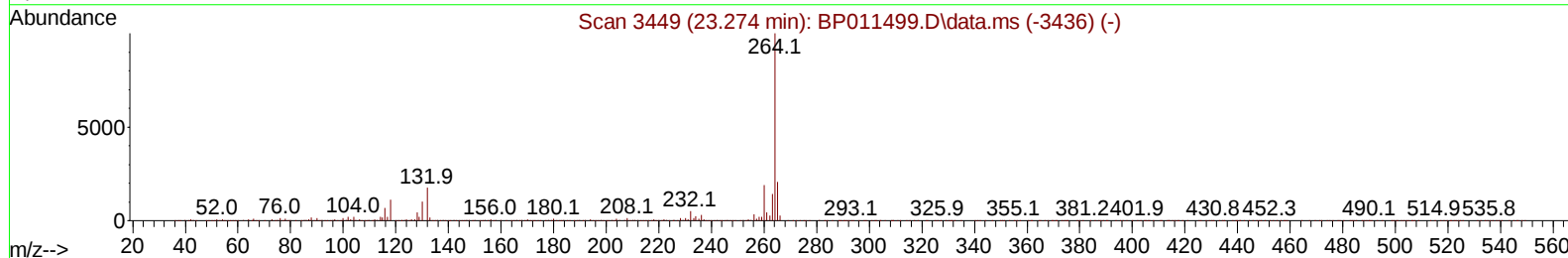
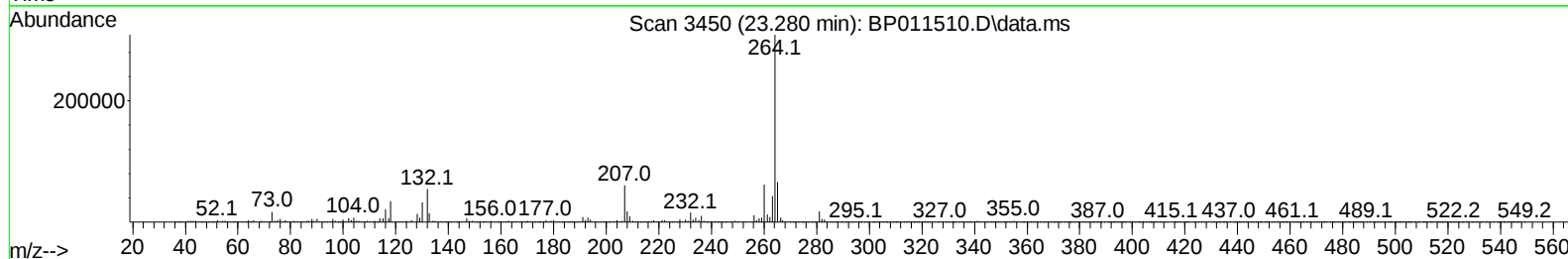
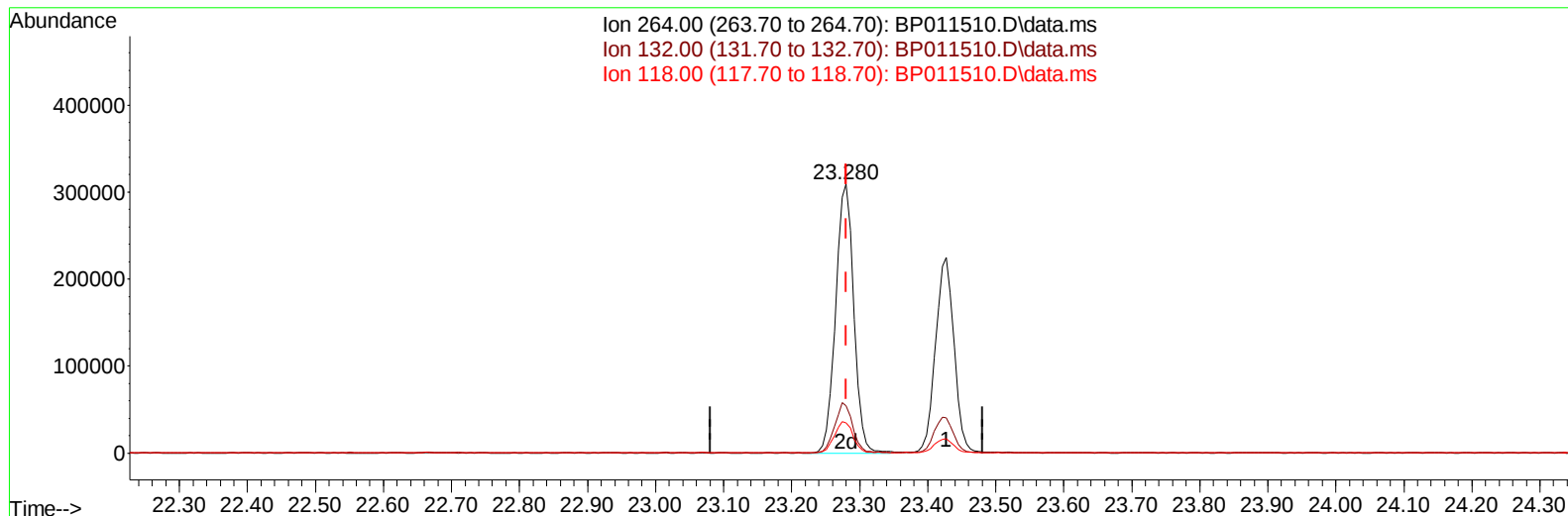
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(92) Benzo(a)pyrene-d12 (S)

23.280min (-0.000) 27.90 ng/ul m

response 574094

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	17.60	17.58
118.00	11.10	11.27
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	7.558	152	98736	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	405536	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.228	164	225357	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	439956	20.000	ng/ul	0.00
79) Chrysene-d12	21.133	240	411879	20.000	ng/ul	0.00
88) Perylene-d12	23.427	264	423033	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	11597	3.645	ng/uL	0.00
4) Pyridine-d5	3.487	84	73847	9.065	ng/ul	0.00
7) Phenol-d5	6.740	99	184474	20.331	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	119275	20.234	ng/ul	0.00
11) 2-Chlorophenol-d4	7.093	132	145812	21.850	ng/ul	0.00
15) 4-Methylphenol-d8	8.281	113	151839	20.908	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	71030	24.329	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	73480	24.730	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	126572	23.077	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	196234	21.493	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	425969	26.725	ng/ul	0.00
49) Acenaphthylene-d8	13.922	160	471789	26.632	ng/ul	0.00
54) 4-Nitrophenol-d4	14.463	143	58734	19.126	ng/ul	0.00
60) Fluorene-d10	15.234	176	348936	26.022	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	40127	16.879	ng/ul	0.00
73) Anthracene-d10	17.104	188	541405	27.786	ng/ul	0.00
81) Pyrene-d10	19.363	212	593344	28.473	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.280	264	574094m	27.899	ng/ul	0.00

Target Compounds Qvalue

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

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