

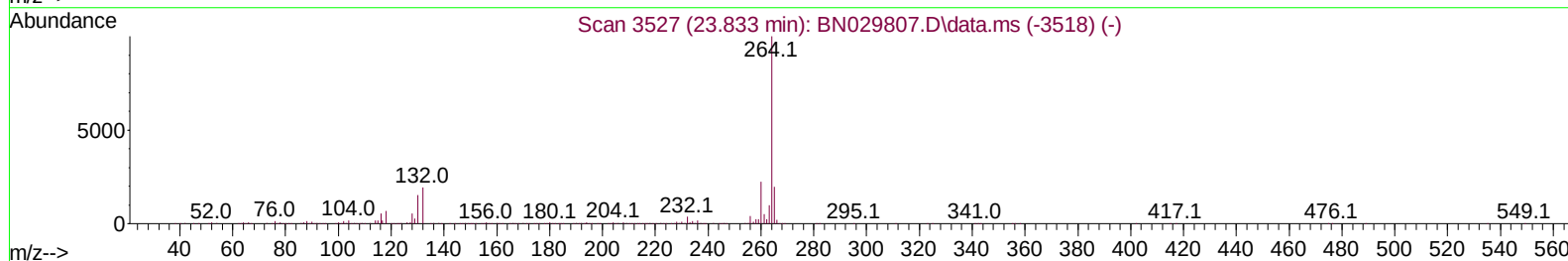
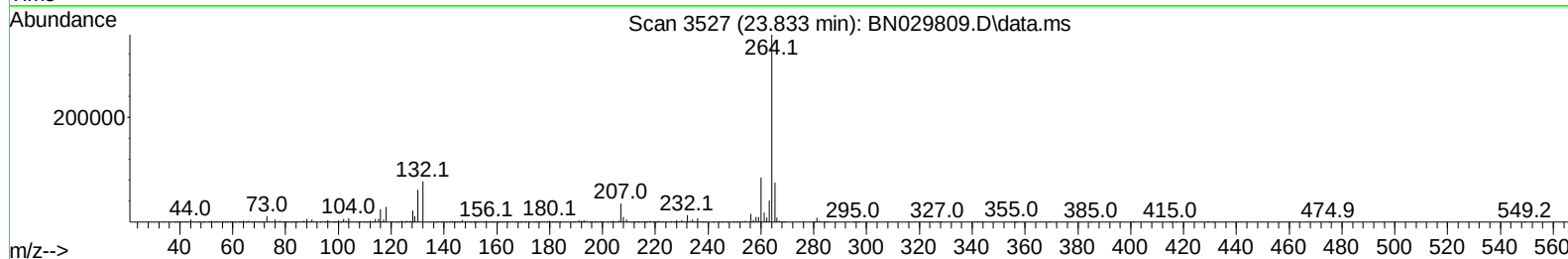
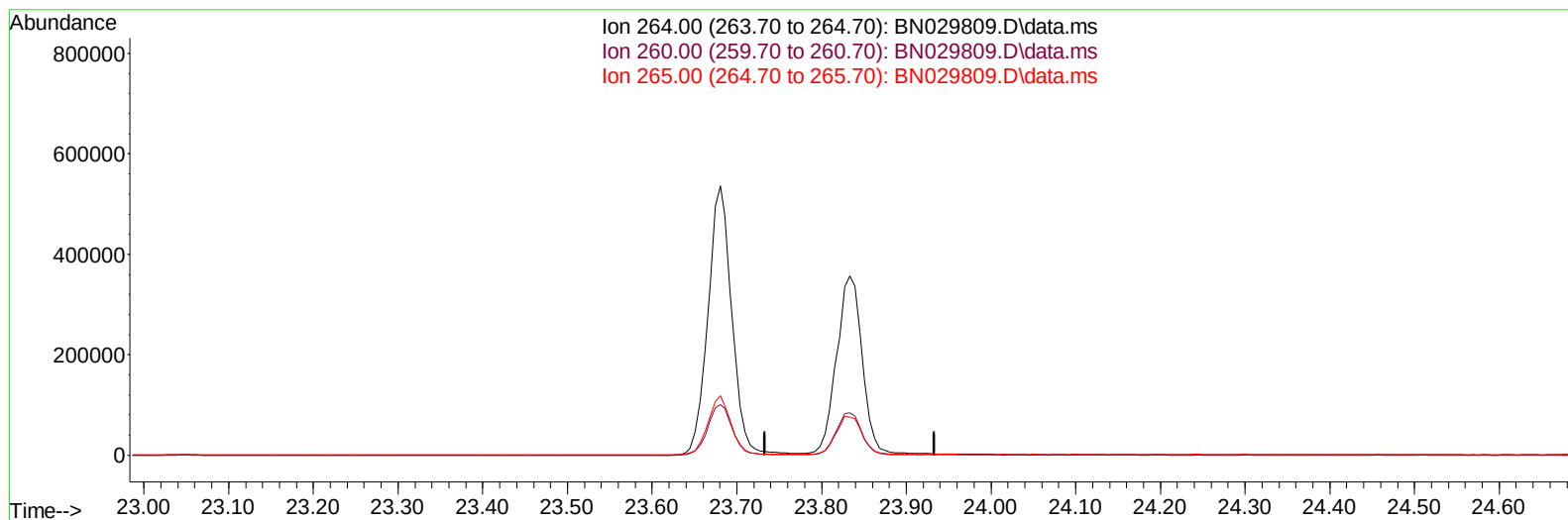
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029809.D
Acq On : 15 Feb 2024 16:05
Operator : MA/JU
Sample : PB158893BL
Misc :
ALS Vial : 89 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK893

Manual IntegrationsAPPROVED

Quant Time: Feb 15 16:32:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 13 22:50:45 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 02/16/2024
Supervised By :mohammad ahmed 02/17/2024



TIC: BN029809.D\data.ms

(88) Perylene-d12 (I)

23.833min (-23.833) 0.00 ng/ul

response 0

Ion	Exp%	Act%
264.00	100.00	0.00
260.00	23.40	0.00#
265.00	20.20	0.00#
0.00	0.00	0.00

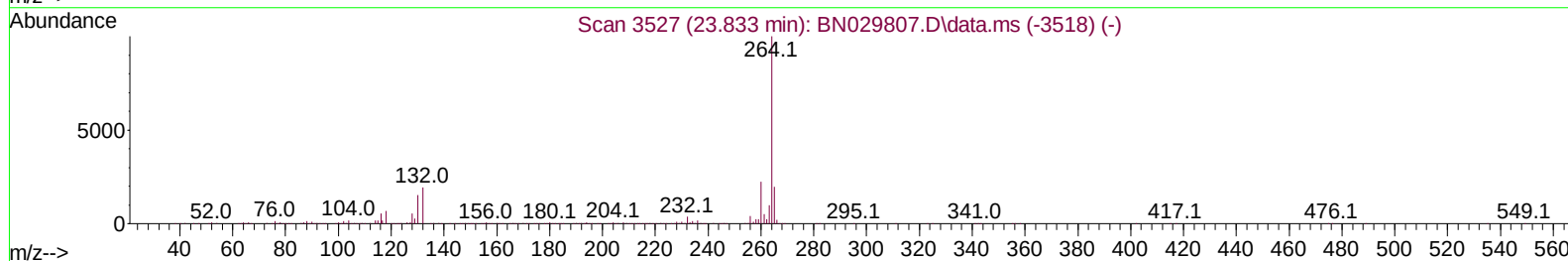
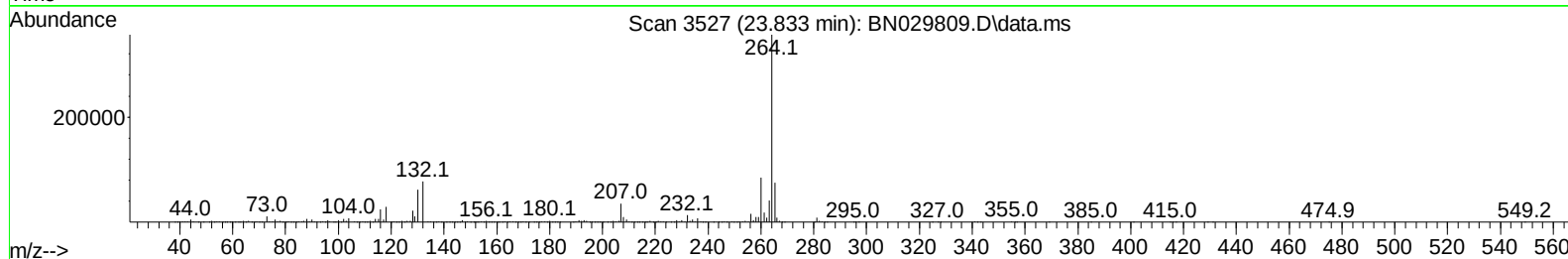
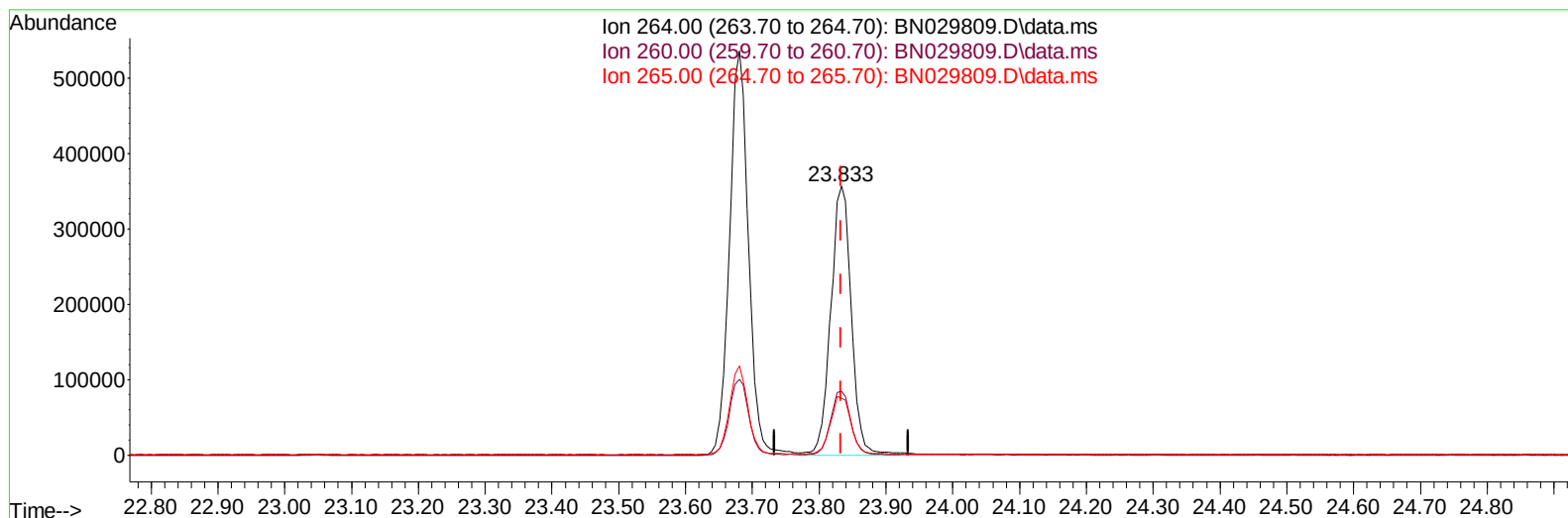
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(88) Perylene-d12 (I)

23.833min (-0.000) 20.00 ng/ul m

response 762217

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	23.40	23.62
265.00	20.20	21.11
0.00	0.00	0.00

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 QLast Update : Tue Feb 13 22:50:45 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	162749	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	665252	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164	352618	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	737853	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	680557	20.000	ng/ul	0.00
88) Perylene-d12	23.833	264	762217m	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	22839	4.616	ng/uL	0.00
4) Pyridine-d5	3.728	84	293217	22.221	ng/ul	0.00
7) Phenol-d5	7.034	99	361666	23.758	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	233193	22.953	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	288303	26.135	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	280758	24.730	ng/ul	0.00
21) Nitrobenzene-d5	9.034	128	137489	26.629	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	138282	30.751	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	239495	26.983	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	374103	24.706	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	725893	27.650	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	836334	26.825	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	108395	21.754	ng/ul	0.00
60) Fluorene-d10	15.504	176	603647	27.029	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	72313	20.113	ng/ul	0.00
73) Anthracene-d10	17.363	188	955172	27.622	ng/ul	0.00
81) Pyrene-d10	19.663	212	1081508	26.481	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	1053920	27.633	ng/ul	0.00

Target Compounds Qvalue

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

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