

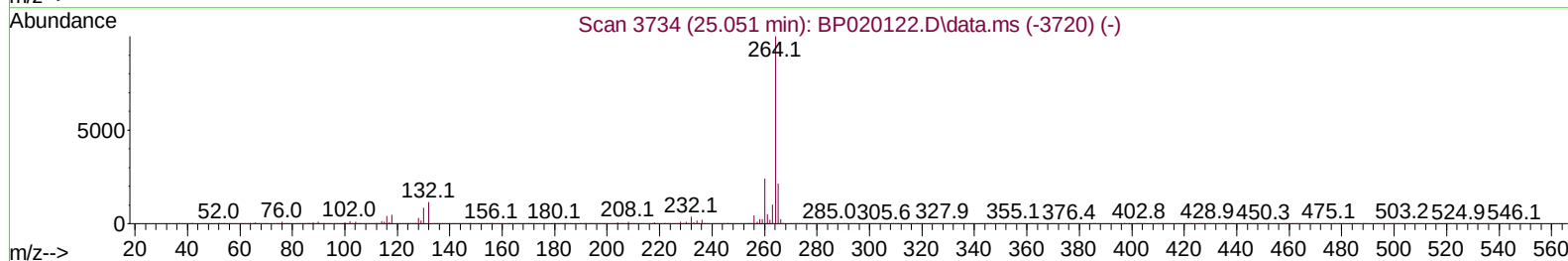
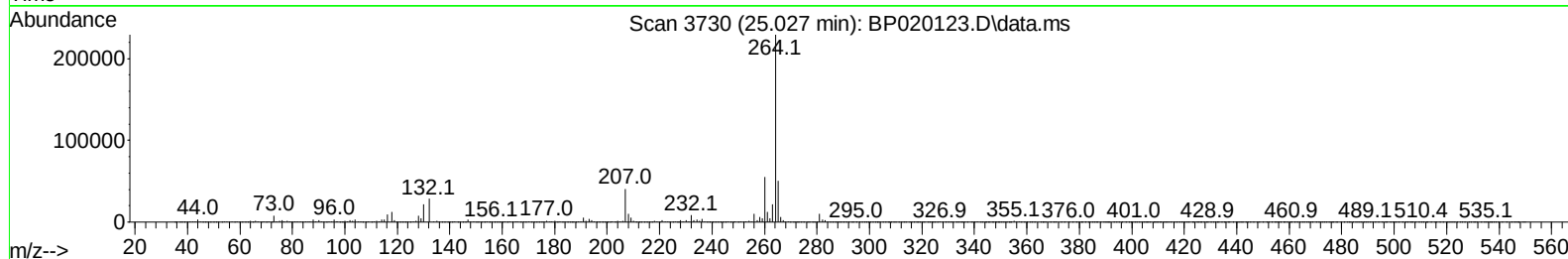
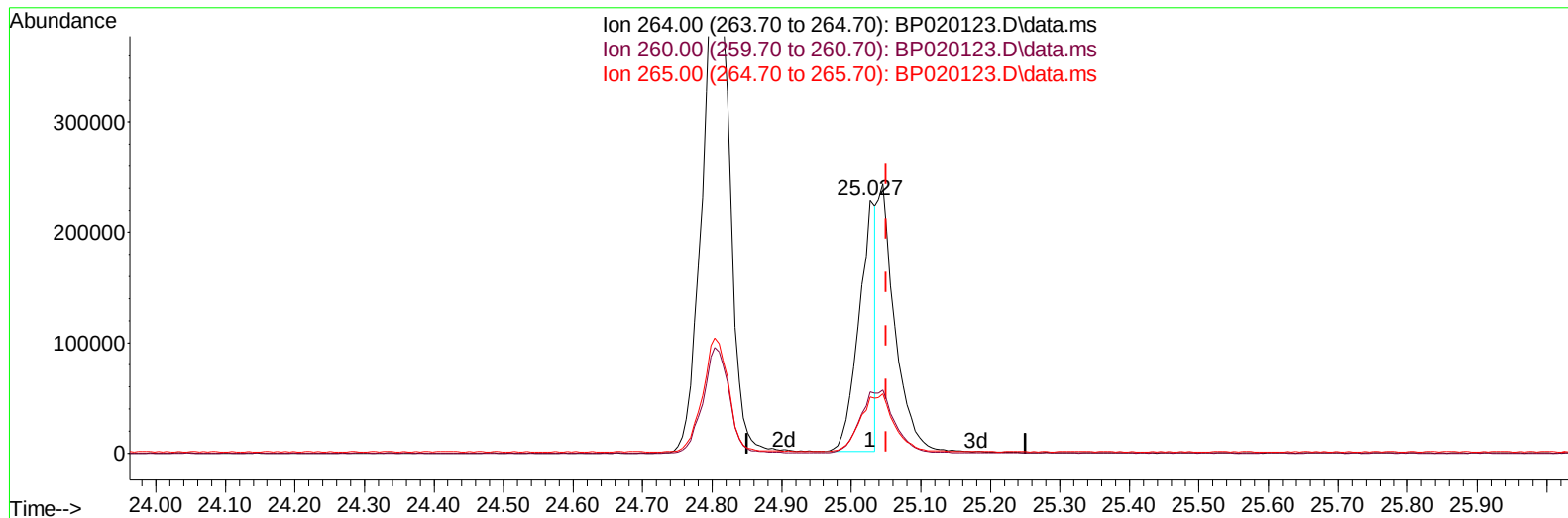
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP043024\
 Data File : BP020123.D
 Acq On : 30 Apr 2024 17:22
 Operator : MA/JU
 Sample : PB160629BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK629

Manual IntegrationsAPPROVED

Quant Time: May 01 00:58:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 01 00:58:38 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024
 Supervised By :mohammad ahmed 05/03/2024



TIC: BP020123.D\data.ms

(88) Perylene-d12 (I)

25.027min (-0.023) 20.00 ng/ul

response 376713

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	23.20	24.20
265.00	21.90	22.14
0.00	0.00	0.00

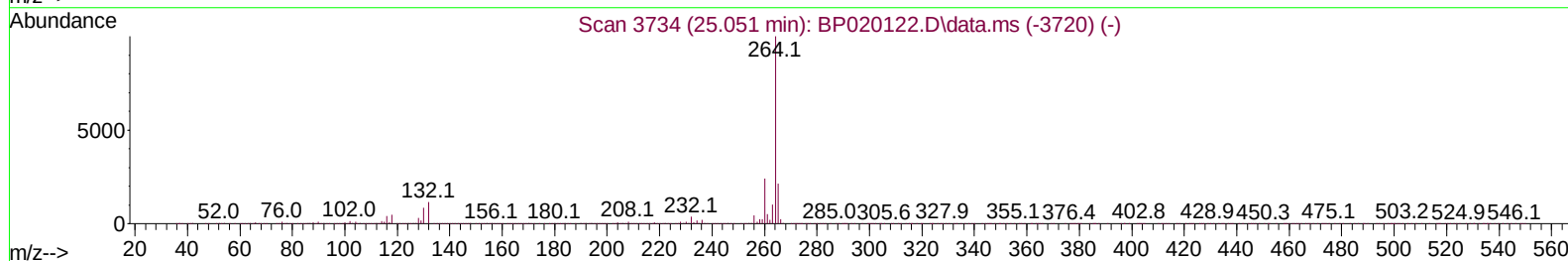
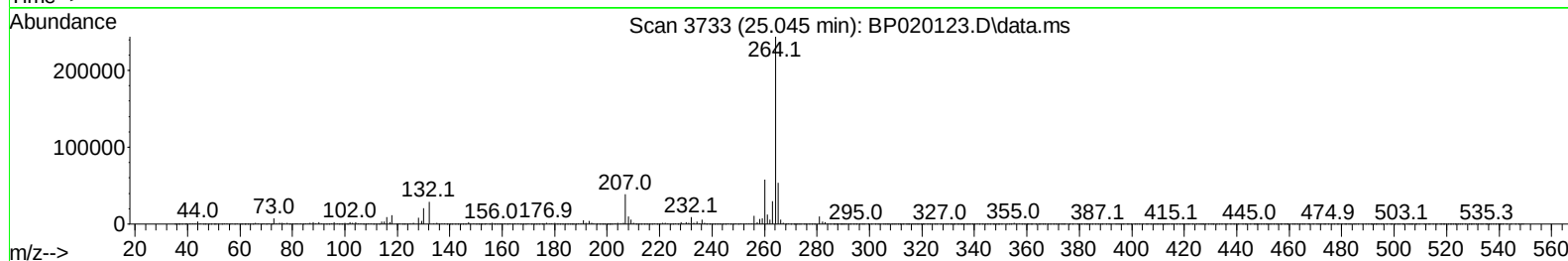
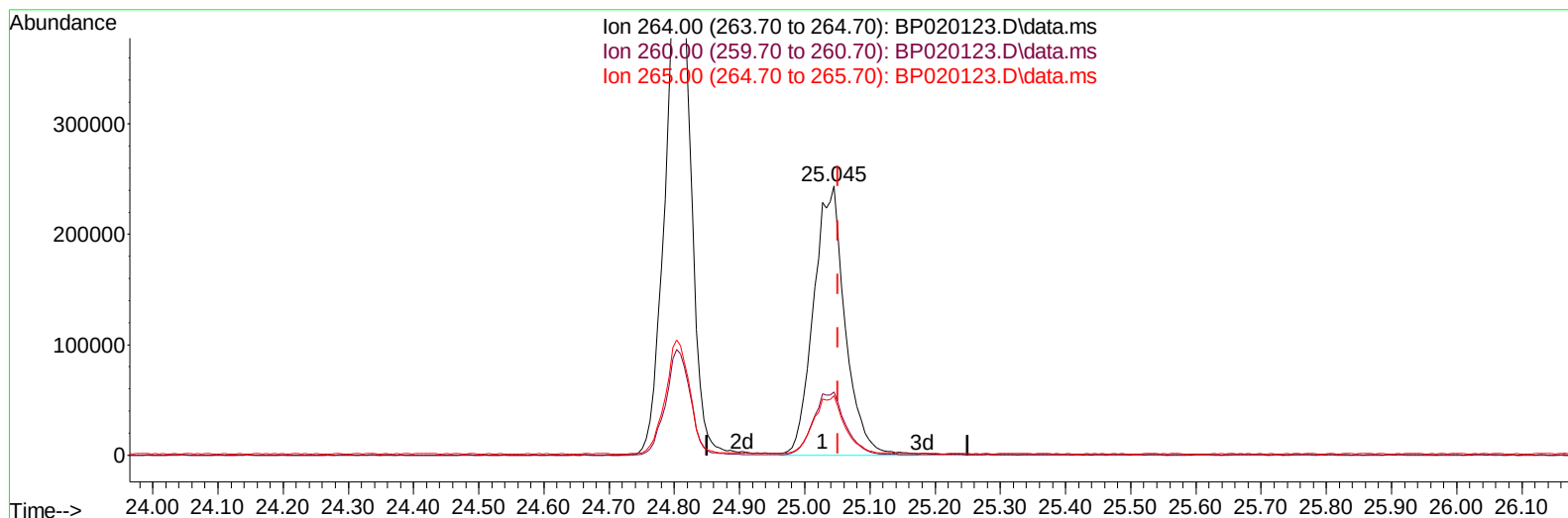
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(88) Perylene-d12 (I)

25.045min (-0.006) 20.00 ng/ul m

response 821532

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	23.20	23.58
265.00	21.90	22.06
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.634	152		95813	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136		408104	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164		316334	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188		754512	20.000	ng/ul	0.00
79) Chrysene-d12	21.657	240		792193	20.000	ng/ul	-0.01
88) Perylene-d12	25.045	264		821532m	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.217	96		12805	5.581	ng/uL	0.00
4) Pyridine-d5	3.623	84		196776	29.113	ng/ul	0.00
7) Phenol-d5	6.822	99		247778	30.497	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67		172258	32.912	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132		181227	31.770	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113		200592	32.429	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128		95754	30.553	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143		107834	31.004	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165		188327	28.772	ng/ul	0.01
31) 4-Chloroaniline-d4	10.569	131		276736	32.421	ng/ul	0.01
46) Dimethylphthalate-d6	13.734	166		771847	34.162	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160		785692	30.510	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143		107088	29.159	ng/ul	0.00
60) Fluorene-d10	15.345	176		637438	32.435	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200		124570	25.821	ng/ul	0.00
73) Anthracene-d10	17.263	188		1076690	32.140	ng/ul	0.00
81) Pyrene-d10	19.663	212		1356560	32.187	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.804	264		1259036	31.409	ng/ul	0.00

Target Compounds Qvalue

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

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