

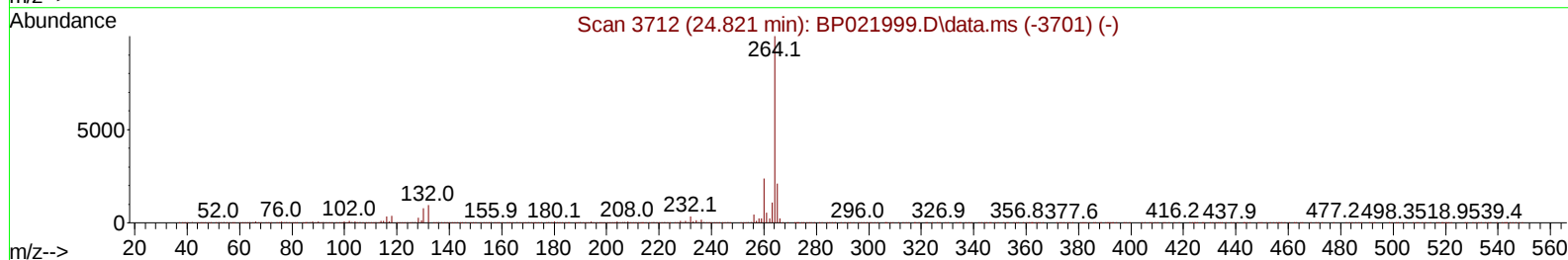
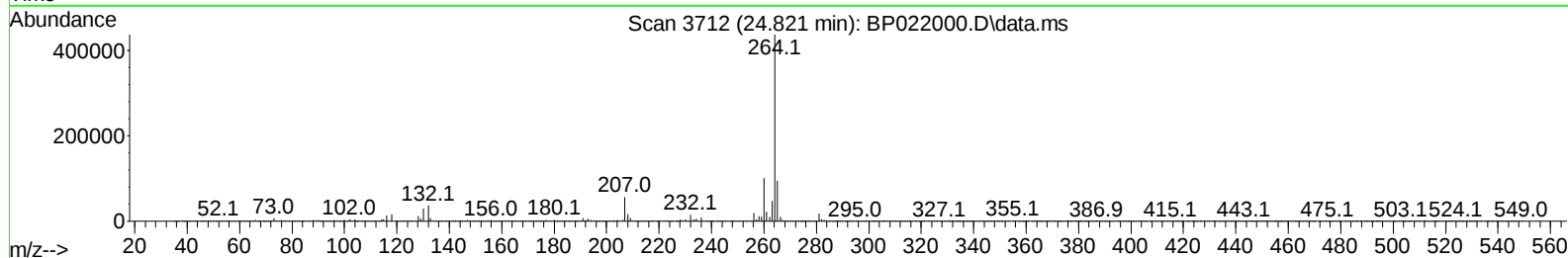
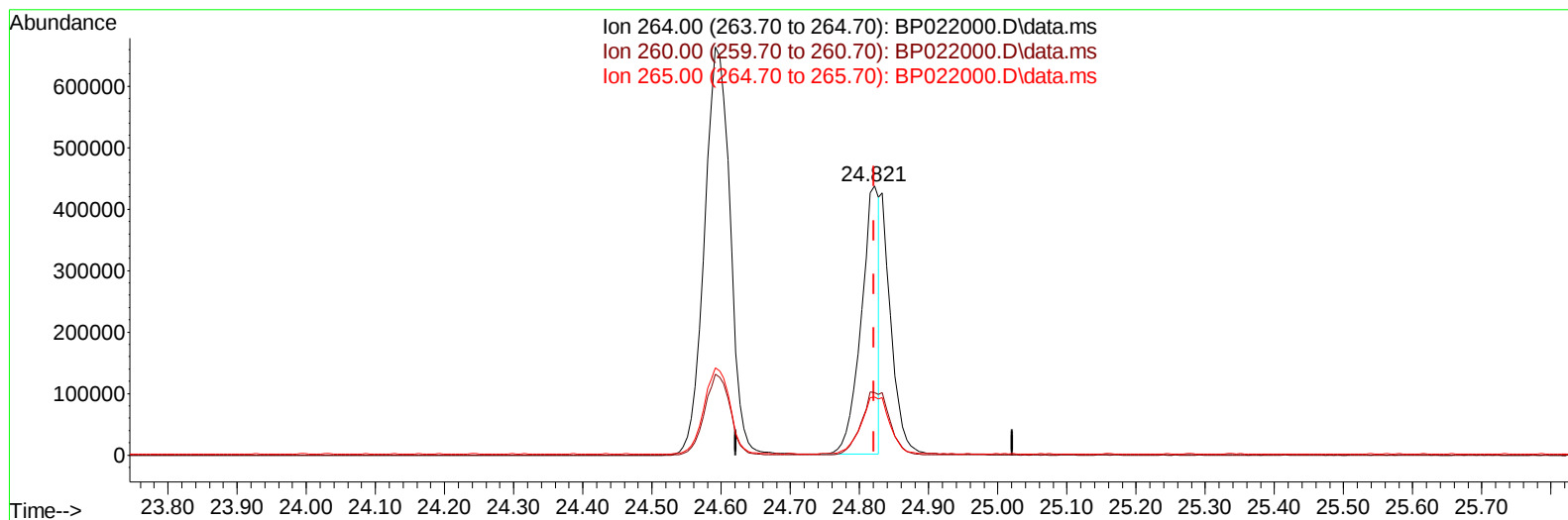
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022000.D
 Acq On : 24 Sep 2024 18:20
 Operator : RC/JU
 Sample : PB163502BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK502

Manual IntegrationsAPPROVED

Quant Time: Sep 24 18:48:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/25/2024
 Supervised By :mohammad ahmed 09/28/2024



TIC: BP022000.D\data.ms

(88) Perylene-d12 (I)

24.821min (-0.000) 20.00 ng/u1

response 788471

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	23.60	23.22
265.00	21.60	21.65
0.00	0.00	0.00

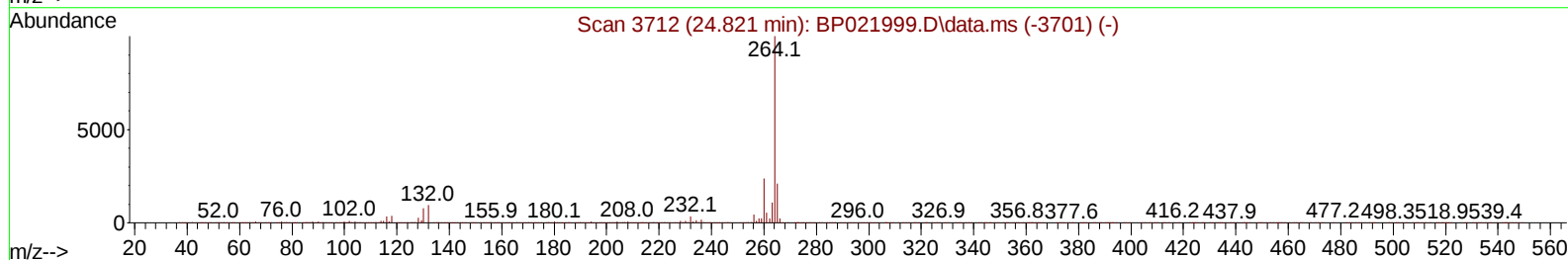
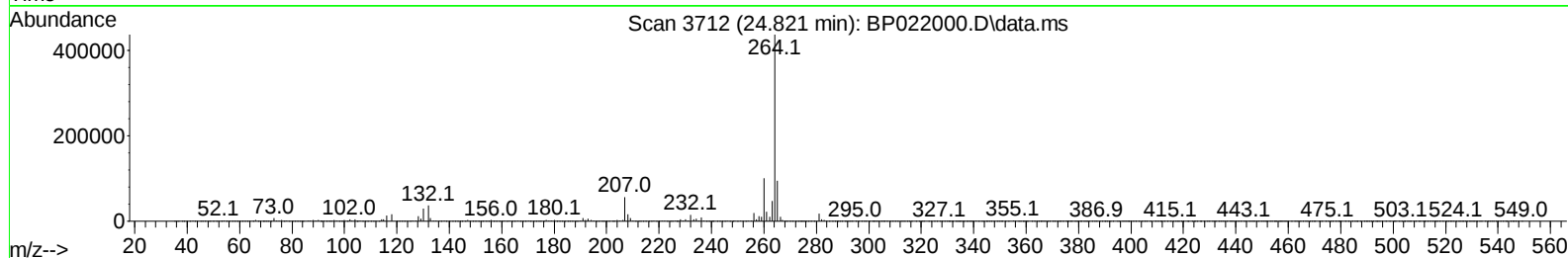
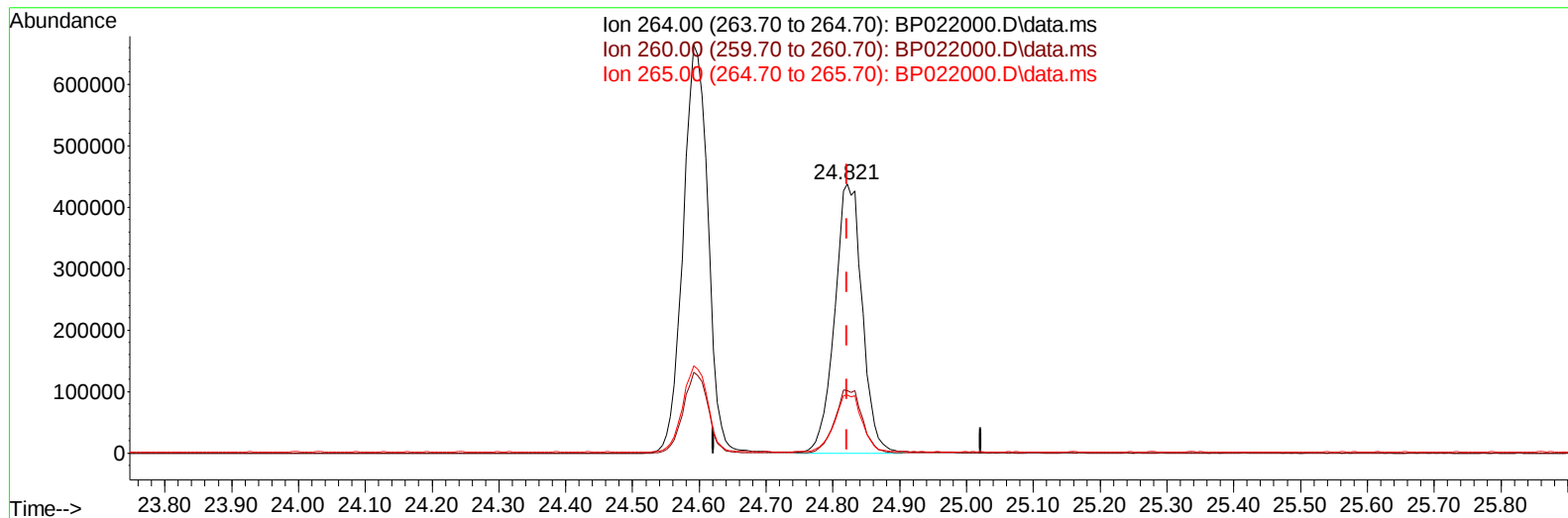
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(88) Perylene-d12 (I)

24.821min (-0.000) 20.00 ng/ul m

response 1248775

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	23.60	23.22
265.00	21.60	21.65
0.00	0.00	0.00

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Reviewed By :Yogesh Patel 09/25/2024
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Quant Time: Sep 24 18:49:28 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	129393	20.000	ng/ul	0.00
20) Naphthalene-d8	10.475	136	471345	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.322	164	364452	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	909981	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1078364	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264	1248775m	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	14674	4.442	ng/uL	0.00
4) Pyridine-d5	3.634	84	217050	22.986	ng/ul	0.00
7) Phenol-d5	6.887	99	249429	23.566	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	156269	23.977	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	188716	24.643	ng/ul	0.00
15) 4-Methylphenol-d8	8.404	113	193631	23.087	ng/ul	0.00
21) Nitrobenzene-d5	8.851	128	90951	25.781	ng/ul	0.00
24) 2-Nitrophenol-d4	9.563	143	103708	25.927	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.104	165	204716	23.566	ng/ul	0.01
31) 4-Chloroaniline-d4	10.598	131	260301	25.114	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	794992	28.219	ng/ul	0.01
49) Acenaphthylene-d8	14.016	160	770158	25.820	ng/ul	0.01
54) 4-Nitrophenol-d4	14.528	143	112582	23.362	ng/ul	0.00
60) Fluorene-d10	15.328	176	649287	26.054	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	130813	23.154	ng/ul	0.01
73) Anthracene-d10	17.222	188	1167110	27.473	ng/ul	0.00
81) Pyrene-d10	19.592	212	1497299	26.801	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.592	264	1717059	26.142	ng/ul	-0.02

Target Compounds Qvalue

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

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