

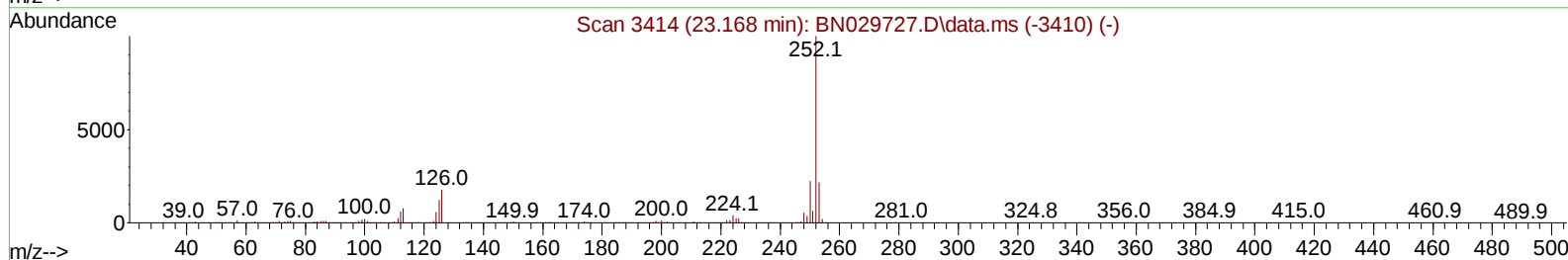
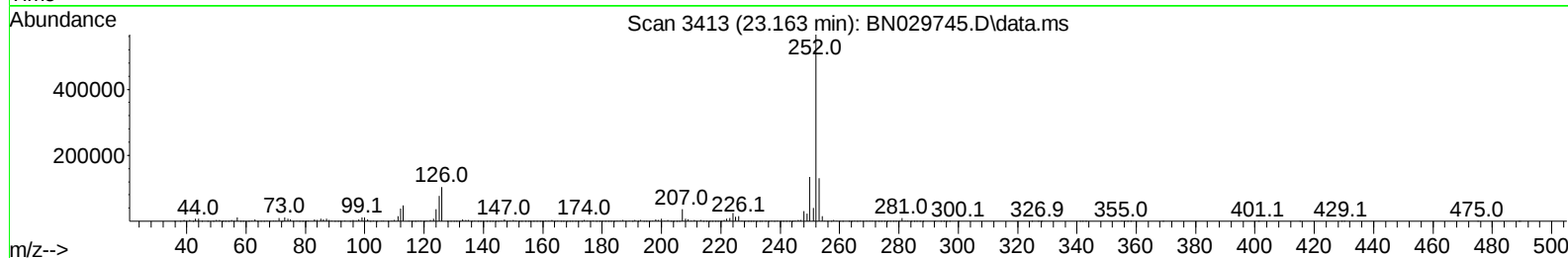
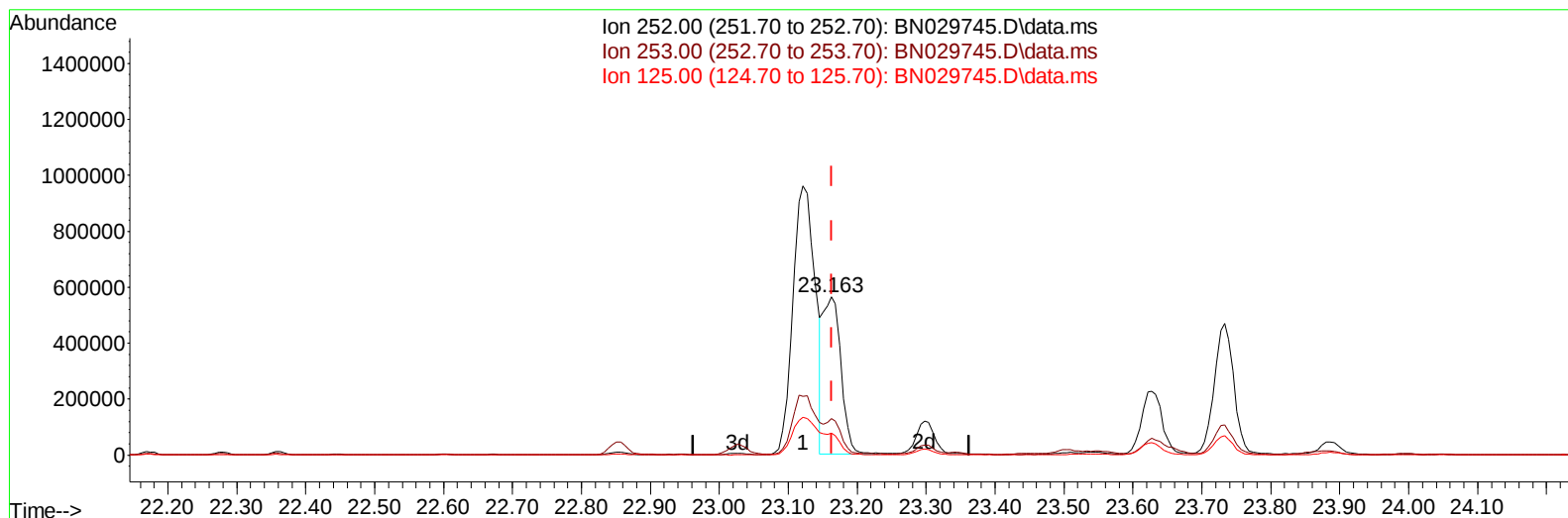
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029745.D
Acq On : 13 Feb 2024 23:04
Operator : MA/JU
Sample : P1352-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0FN3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/15/2024
Supervised By :mohammad ahmed 02/17/2024

Quant Time: Feb 14 00:12:17 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



TIC: BN029745.D\data.ms

(91) Benzo(k)fluoranthene

23.163min (-0.000) 20.66 ng/u1 m

response 1021304

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.80	23.00
125.00	14.00	13.68
0.00	0.00	0.00

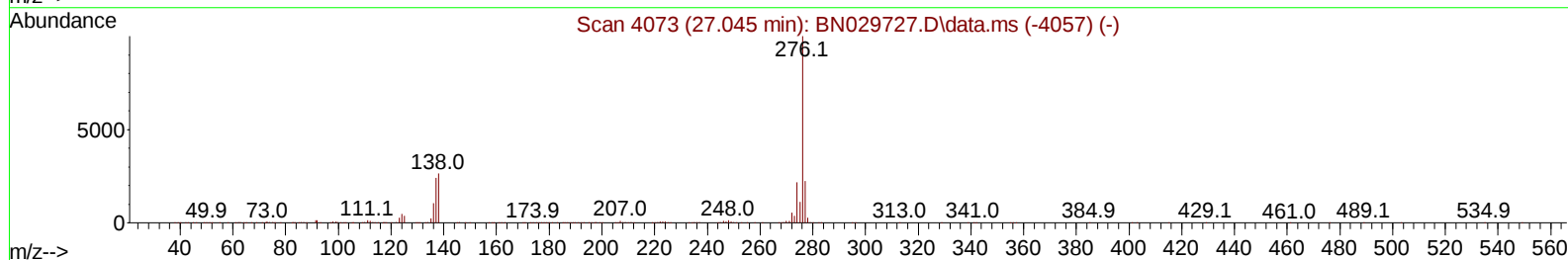
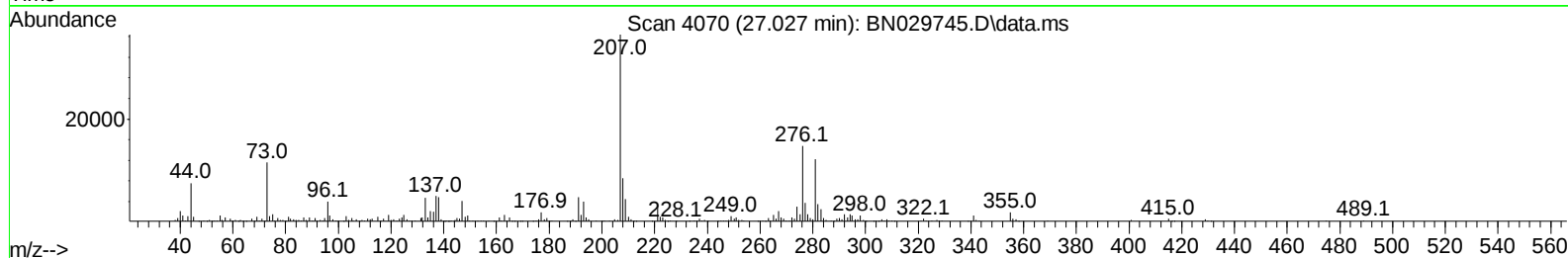
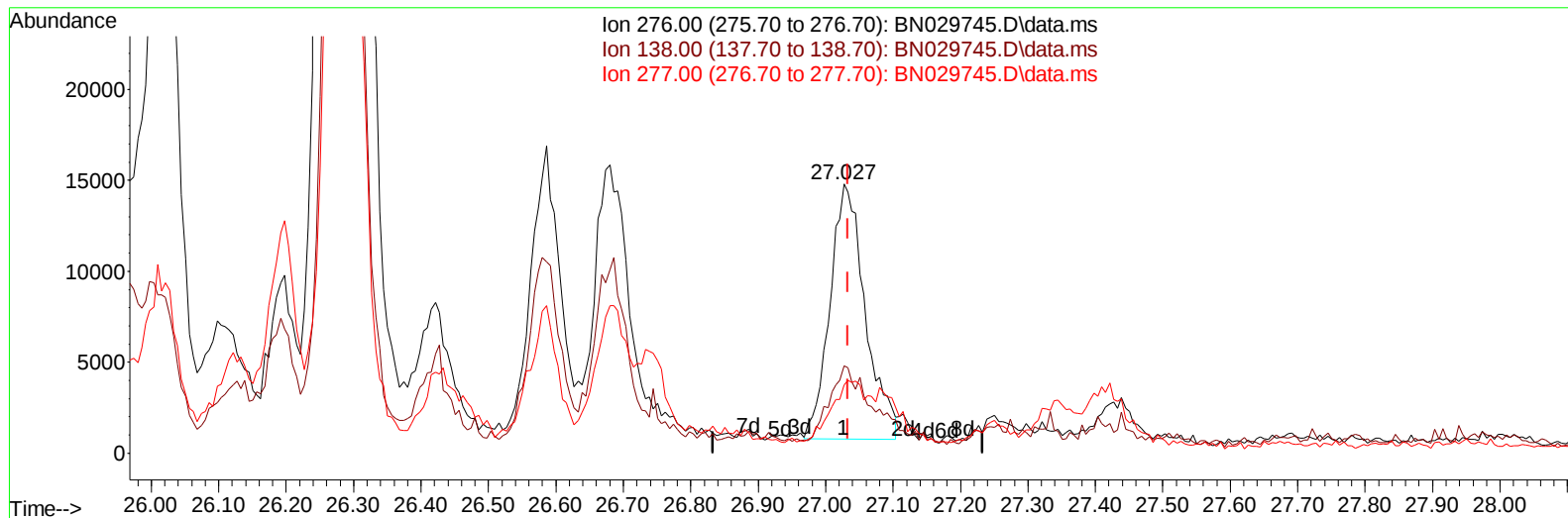
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TIC: BN029745.D\data.ms

(96) Benzo(g,h,i)perylene

27.027min (-0.006) 1.08 ng/u1

response 50918

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	29.90	32.42
277.00	23.30	25.34
0.00	0.00	0.00

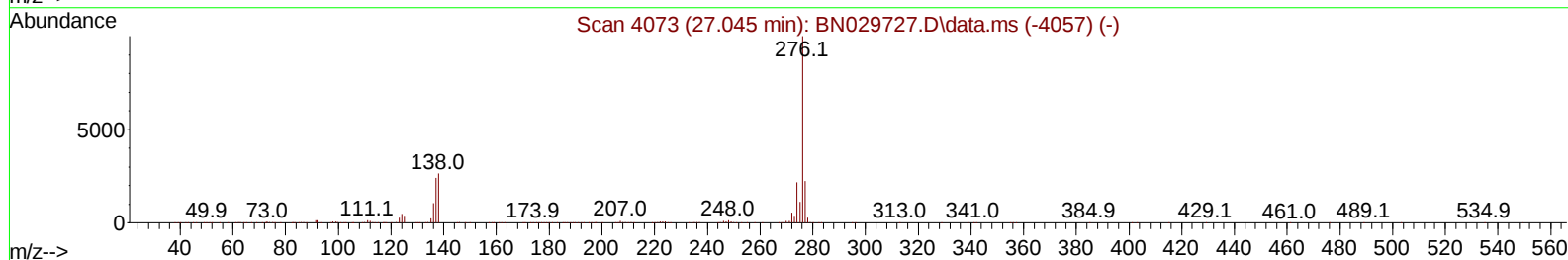
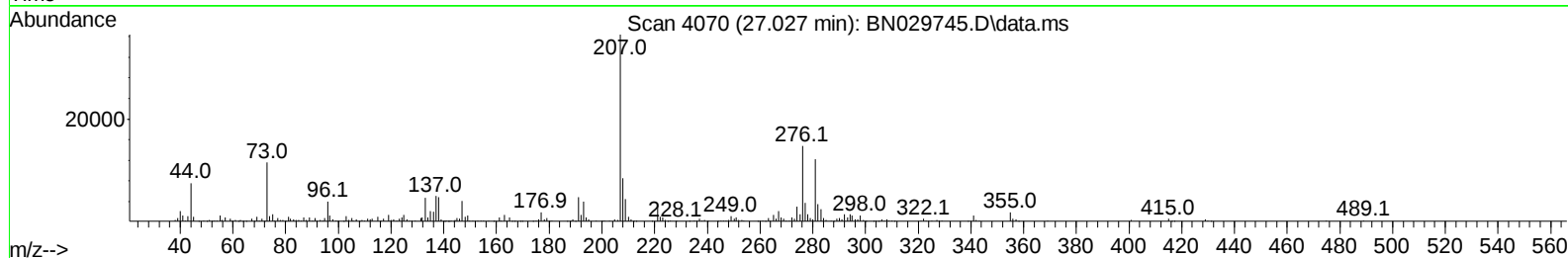
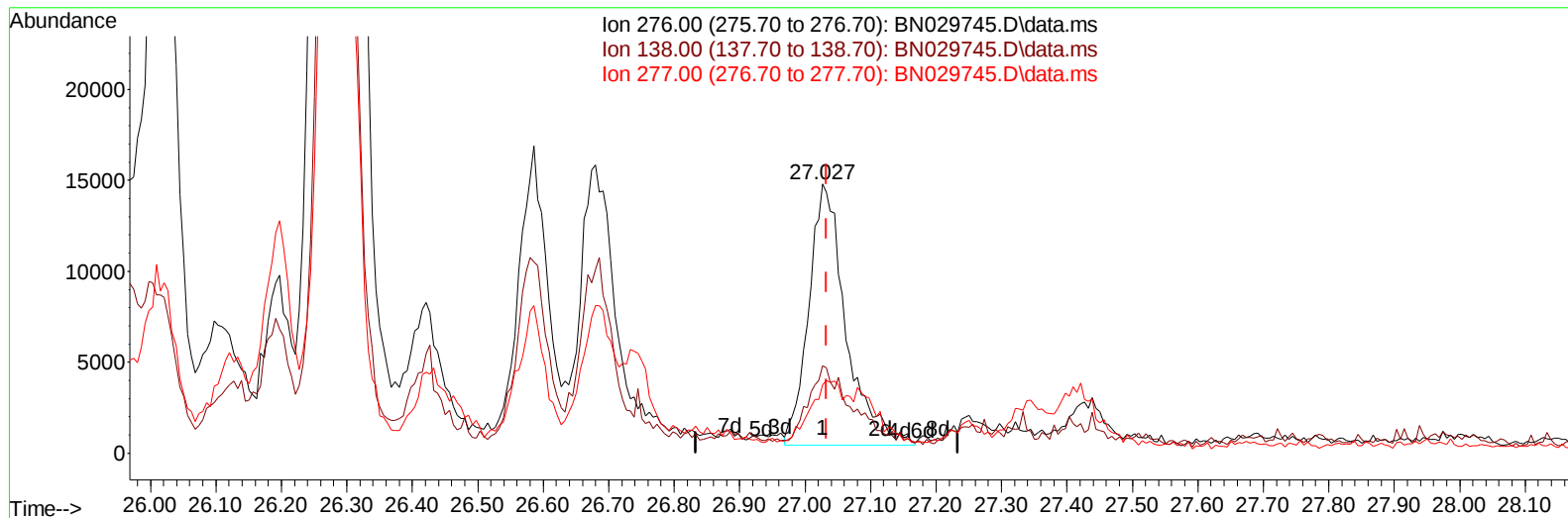
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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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TIC: BN029745.D\data.ms

(96) Benzo(g,h,i)perylene

27.027min (-0.006) 1.20 ng/ul m

response 56674

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	29.90	32.42
277.00	23.30	25.34
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/15/2024
 Supervised By :mohammad ahmed 02/17/2024

Quant Time: Feb 14 00:15:05 2024
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 QLast Update : Tue Feb 13 22:50:45 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.869	152		208669	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136		809395	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164		377870	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188		685296	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240		630098	20.000	ng/ul	0.00
88) Perylene-d12	23.839	264		766843	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.322	96		31516	4.968	ng/uL	0.00
4) Pyridine-d5	3.734	84		235590	13.925	ng/ul	0.00
7) Phenol-d5	7.040	99		473558	24.263	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67		343911	26.402	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132		390663	27.621	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113		248776	17.091	ng/ul	0.00
21) Nitrobenzene-d5	9.040	128		182534	29.057	ng/ul	0.00
24) 2-Nitrophenol-d4	9.751	143		173920	31.789	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165		291537	26.996	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131		386042	20.954	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166		796006	28.295	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160		983686	29.443	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143		80615	15.098	ng/ul	0.00
60) Fluorene-d10	15.510	176		693495	28.977	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200		60233	18.038	ng/ul	0.00
73) Anthracene-d10	17.363	188		869758	27.081	ng/ul	0.00
81) Pyrene-d10	19.663	212		983413	26.007	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264		819074	21.346	ng/ul	0.00
Target Compounds							
						Qvalue	
16) Acetophenone	8.699	105		37269	1.609	ng/ul	96
52) Acenaphthene	14.575	153		28191	1.064	ng/ul	95
72) Phenanthrene	17.310	178		866229	21.852	ng/ul	99
74) Anthracene	17.398	178		155771	3.915	ng/ul	97
77) Carbazole	17.675	167		84468	2.325	ng/ul	97
80) Fluoranthene	19.333	202		3032102	66.012	ng/ul	99
82) Pyrene	19.692	202		2446013	50.337	ng/ul	99
85) Benzo(a)anthracene	21.451	228		1669187	36.061	ng/ul	96
87) Chrysene	21.504	228		1631303	37.606	ng/ul	99
90) Benzo(b)fluoranthene	23.121	252		2143276	44.373	ng/ul	100
91) Benzo(k)fluoranthene	23.163	252		1021304m	20.658	ng/ul	
93) Benzo(a)pyrene	23.733	252		911831	19.685	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.280	276		707700	12.547	ng/ul	98
95) Dibenzo(a,h)anthracene	26.304	278		239106	5.192	ng/ul	95
96) Benzo(g,h,i)perylene	27.027	276		56674m	1.201	ng/ul	

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

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