

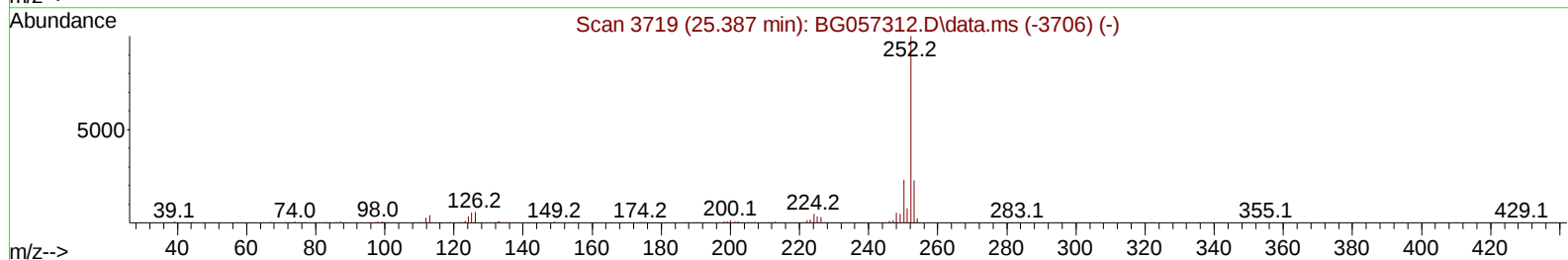
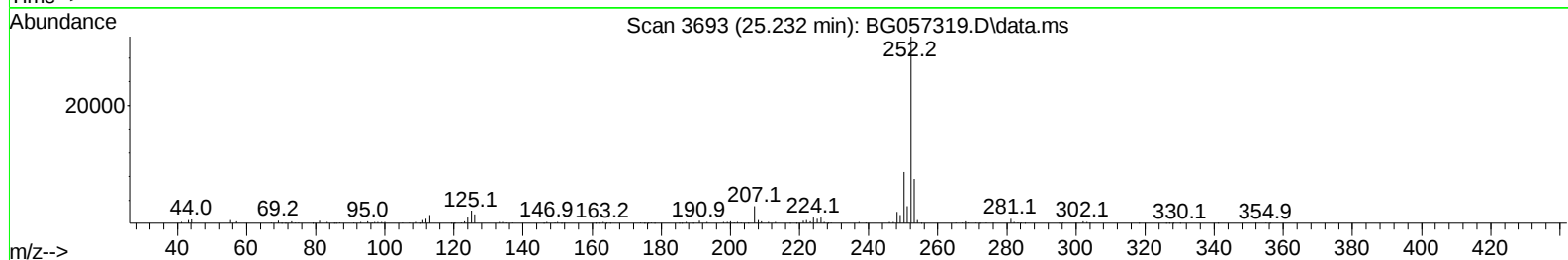
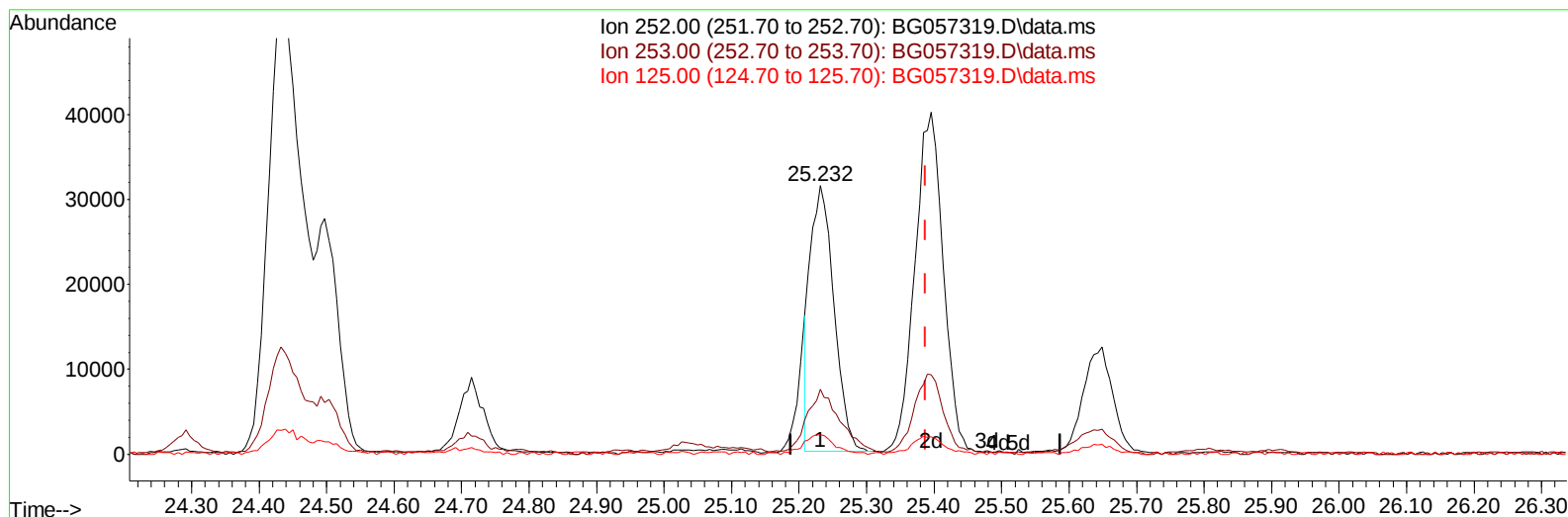
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050523\
Data File : BG057319.D
Acq On : 5 May 2023 15:39
Operator : CG/JU
Sample : 02479-13DL 20X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW976DL

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/08/2023
Supervised By :Jagrut Upadhyay 05/08/2023

Quant Time: May 06 00:52:22 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 05 11:20:08 2023
Response via : Initial Calibration



TIC: BG057319.D\data.ms

(93) Benzo(a)pyrene (C)

25.232min (-0.156) 9.83 ng/u1

response 76638

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.50	24.17
125.00	6.40	7.20
0.00	0.00	0.00

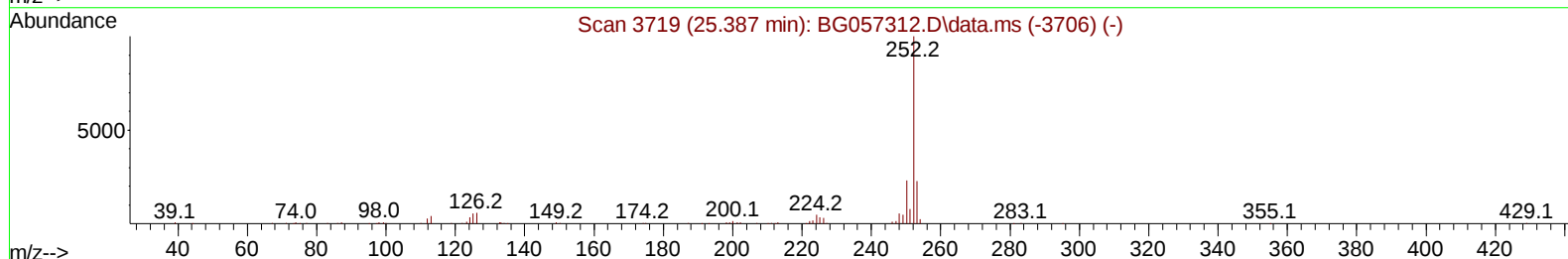
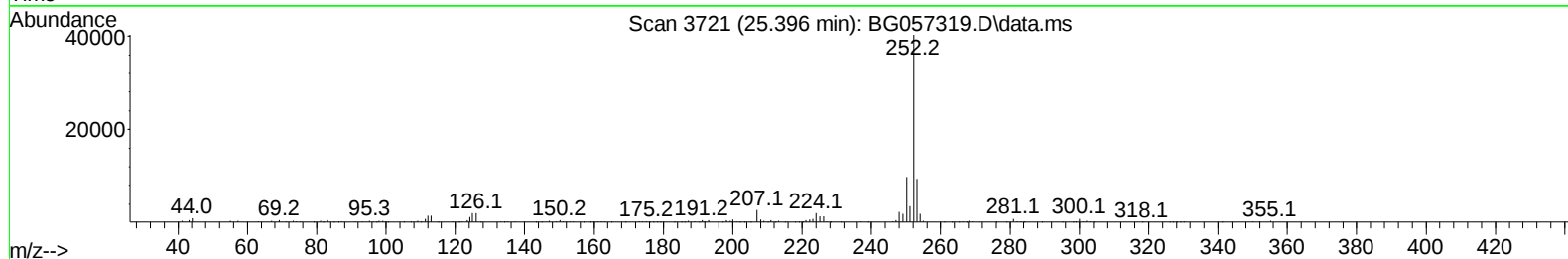
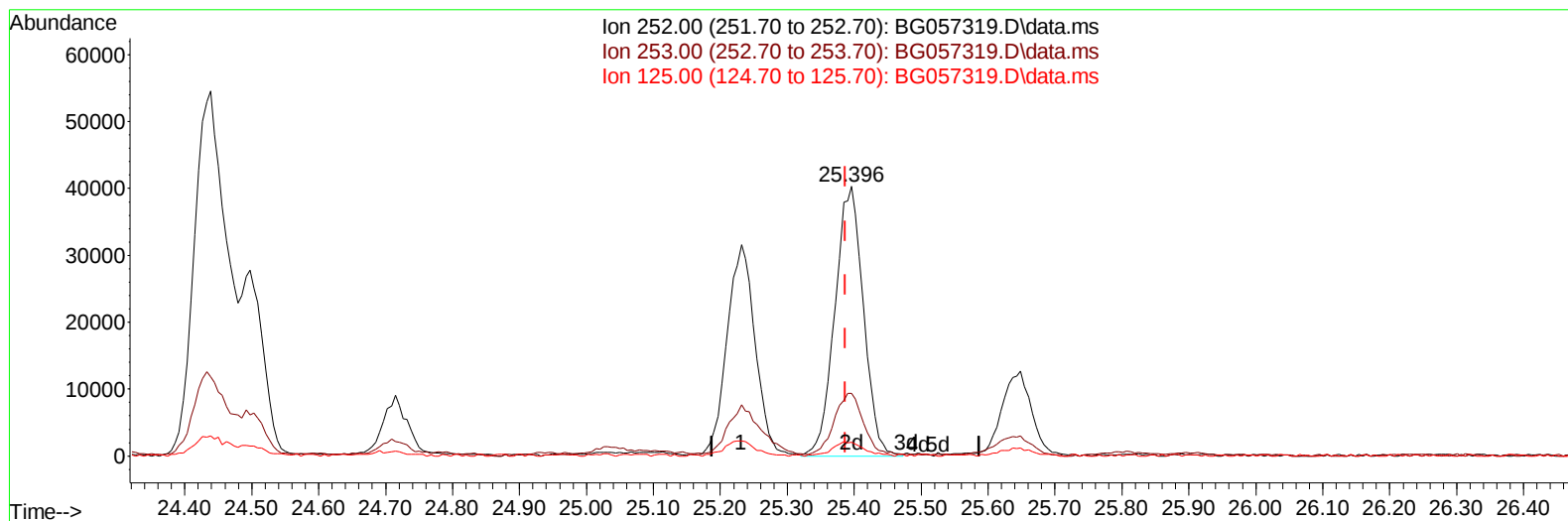
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(93) Benzo(a)pyrene (C)

25.396min (+ 0.009) 15.28 ng/ul m

response 119054

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.50	23.21
125.00	6.40	5.07#
0.00	0.00	0.00

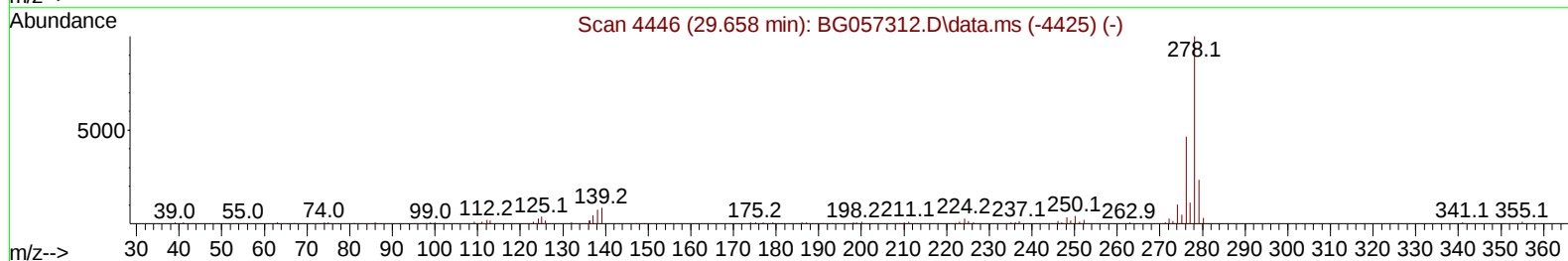
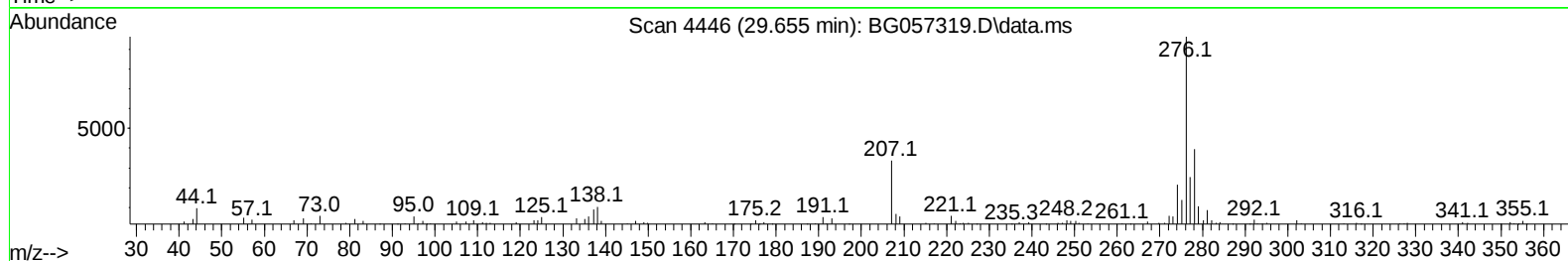
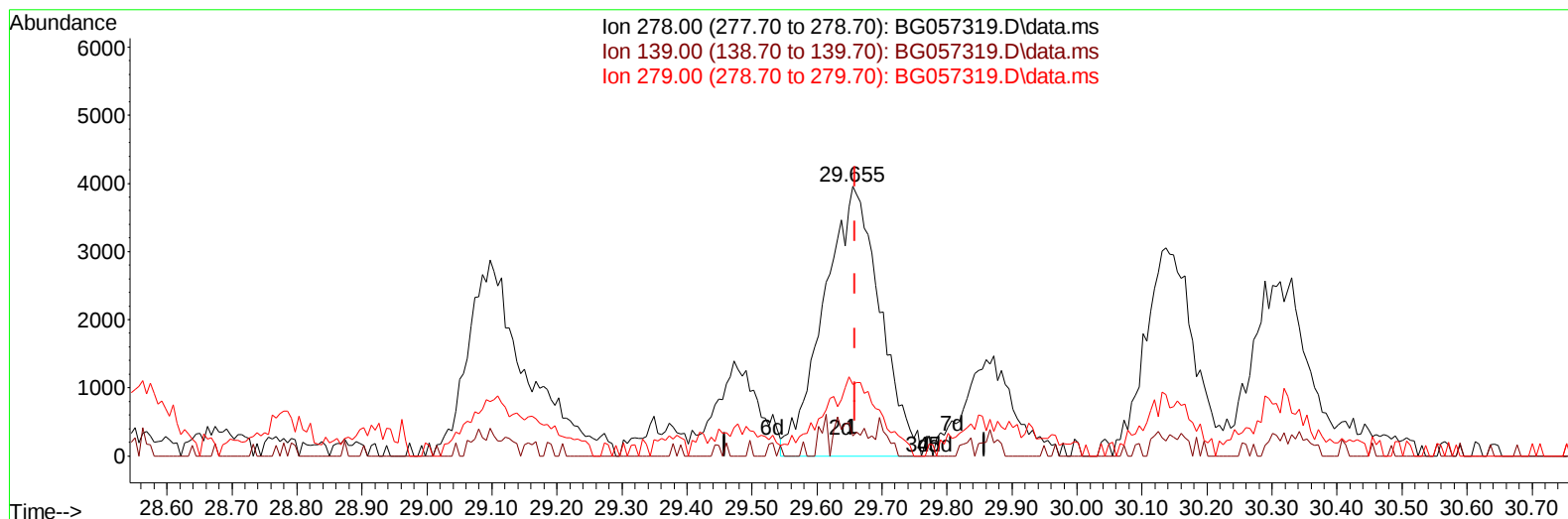
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TIC: BG057319.D\data.ms

(95) Dibenzo(a,h)anthracene

29.655min (-0.003) 2.94 ng/ul m

response 23755

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	7.20	7.53
279.00	22.90	26.45
0.00	0.00	0.00

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Quant Time: May 06 01:45:14 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.372	152		17278	20.000	ng/ul	0.00
20) Naphthalene-d8	11.215	136		78728	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.987	164		62398	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.730	188		142704	20.000	ng/ul	0.00
79) Chrysene-d12	22.042	240		118193	20.000	ng/ul	0.00
88) Perylene-d12	25.560	264		130344	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	0.000	96		0	0.000	ng/uL	
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.520	99		661	0.405	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.679	67		450	0.468	ng/ul	0.00
11) 2-Chlorophenol-d4	7.902	132		569	0.528	ng/ul	0.00
15) 4-Methylphenol-d8	9.089	113		541	0.436	ng/ul	0.01
21) Nitrobenzene-d5	9.547	128		269	0.429	ng/ul	0.00
24) 2-Nitrophenol-d4	0.000	143		0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	10.839	165		745	0.528	ng/ul	0.01
31) 4-Chloroaniline-d4	0.000	131		0	0.000	ng/ul	
46) Dimethylphthalate-d6	14.370	166		3429	0.694	ng/ul	0.00
49) Acenaphthylene-d8	14.687	160		3867	0.696	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143		0	0.000	ng/ul	
60) Fluorene-d10	15.973	176		3142	0.682	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200		0	0.000	ng/ul	
73) Anthracene-d10	17.824	188		5198	0.798	ng/ul	0.00
81) Pyrene-d10	20.091	212		5564	0.793	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.314	264		5352	0.806	ng/ul	0.00
Target Compounds							
						Qvalue	
52) Acenaphthene	15.051	153		7934	1.972	ng/ul	97
56) Dibenzofuran	15.380	168		6458	1.112	ng/ul	93
61) Fluorene	16.026	166		9668	1.968	ng/ul	97
72) Phenanthrene	17.771	178		171963	23.140	ng/ul	99
74) Anthracene	17.859	178		40514	5.427	ng/ul	97
77) Carbazole	18.123	167		13656	2.189	ng/ul	96
80) Fluoranthene	19.762	202		338254	40.589	ng/ul	97
82) Pyrene	20.121	202		282410	33.475	ng/ul#	98
85) Benzo(a)anthracene	22.018	228		141851	16.969	ng/ul	97
87) Chrysene	22.095	228		137818	17.475	ng/ul	98
90) Benzo(b)fluoranthene	24.438	252		184587	20.450	ng/ul	98
91) Benzo(k)fluoranthene	24.497	252		60900	6.935	ng/ul	100
93) Benzo(a)pyrene	25.396	252		119054m	15.278	ng/ul	
94) Indeno(1,2,3-cd)pyrene	29.620	276		83507	8.573	ng/ul	94
95) Dibenzo(a,h)anthracene	29.655	278		23755m	2.941	ng/ul	
96) Benzo(g,h,i)perylene	30.906	276		72492	9.195	ng/ul	100

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

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