

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
 Data File : BG061386.D
 Acq On : 31 May 2024 9:31
 Operator : MA/JU
 Sample : P2570-14
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :

BNA_G

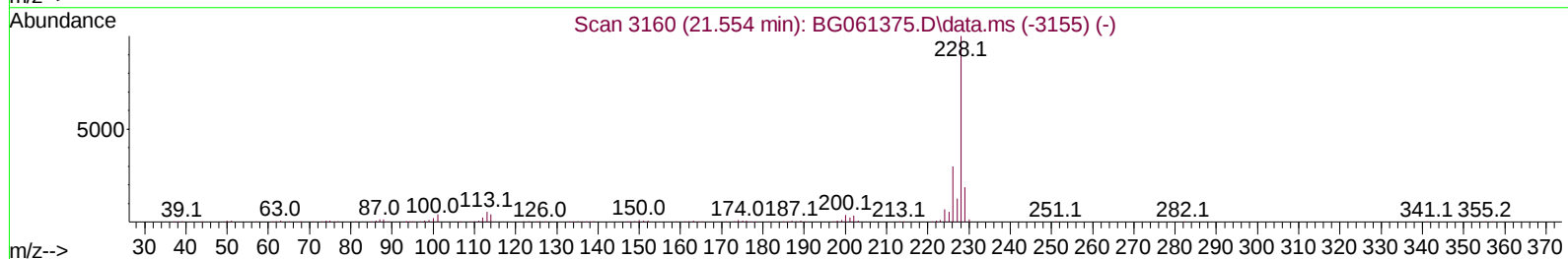
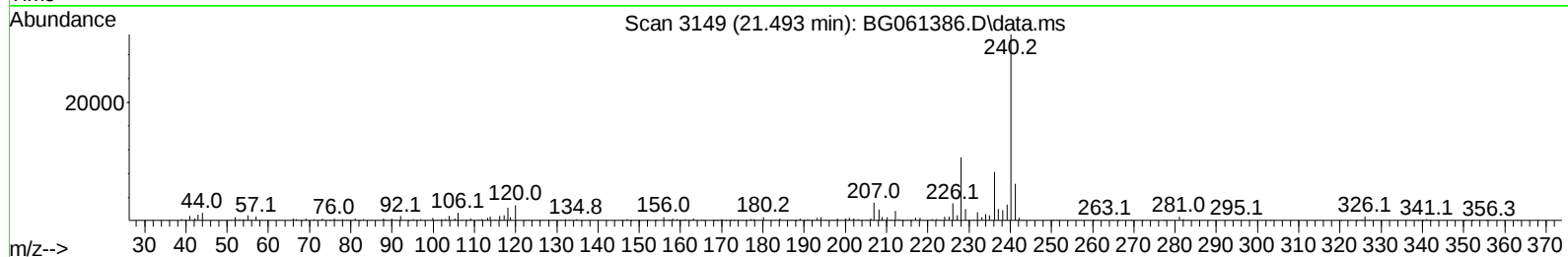
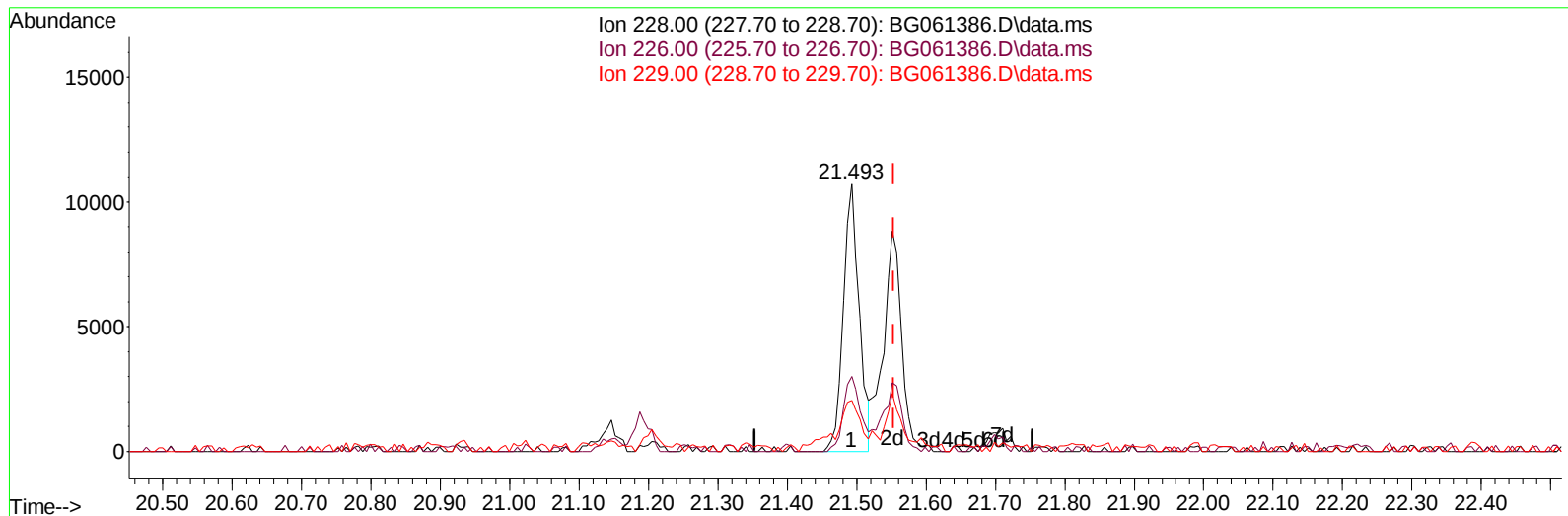
ClientSampleId :

BH5R0

Manual IntegrationsAPPROVED

Quant Time: May 31 10:41:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 30 22:54:48 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/31/2024
 Supervised By :mohammad ahmed 06/01/2024



TIC: BG061386.D\data.ms

(87) Chrysene

21.493min (-0.061) 1.28 ng/ul

response 16880

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.20	27.97
229.00	19.30	19.05
0.00	0.00	0.00

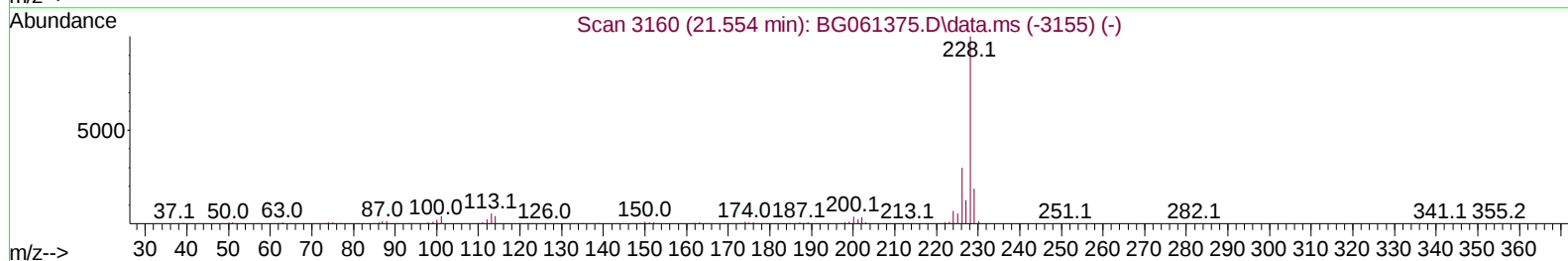
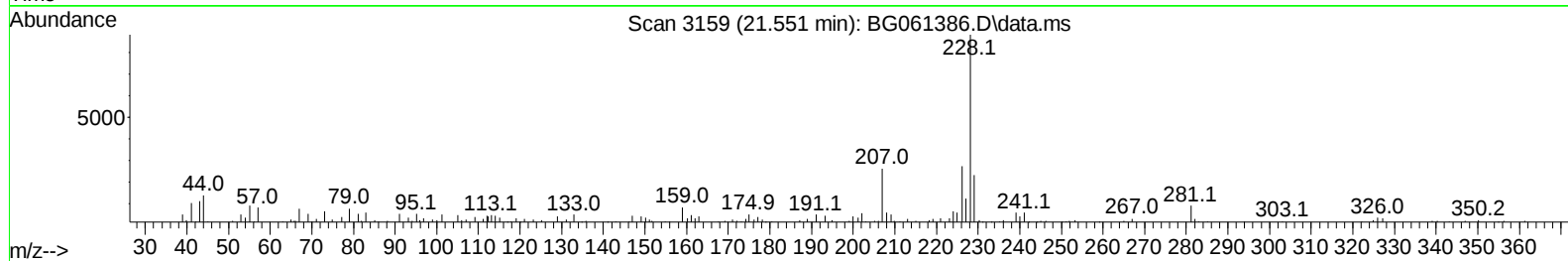
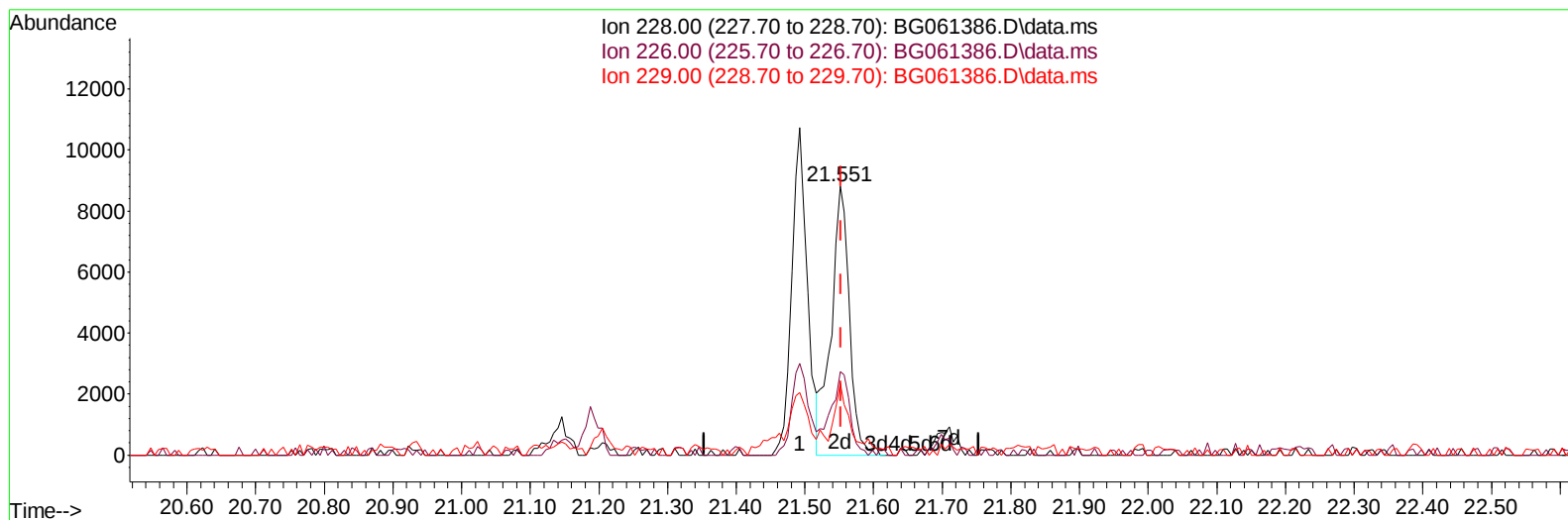
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(87) Chrysene

21.551min (-0.002) 1.23 ng/ul m

response 16278

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.20	31.05
229.00	19.30	26.44#
0.00	0.00	0.00

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Quant Time: May 31 10:45:17 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.873	152	31726	20.000	ng/ul	0.00
20) Naphthalene-d8	10.664	136	118945	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.507	164	87191	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.256	188	196122	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.510	240	207821	20.000	ng/ul	0.00
88) Perylene-d12	24.595	264	234455	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.337	96	2475	4.366	ng/uL	0.00
4) Pyridine-d5	3.743	84	36678	18.186	ng/ul	0.00
7) Phenol-d5	7.057	99	56941	21.855	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.204	67	37071	22.638	ng/ul	0.00
11) 2-Chlorophenol-d4	7.409	132	47526	24.648	ng/ul	0.00
15) 4-Methylphenol-d8	8.590	113	50197	22.595	ng/ul	0.00
21) Nitrobenzene-d5	9.031	128	23924	26.124	ng/ul	0.00
24) 2-Nitrophenol-d4	9.754	143	27208	25.386	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.300	165	54861	26.256	ng/ul	0.00
31) 4-Chloroaniline-d4	10.805	131	63435	23.205	ng/ul	0.00
46) Dimethylphthalate-d6	13.913	166	178922	26.459	ng/ul	0.00
49) Acenaphthylene-d8	14.201	160	199318	28.351	ng/ul	0.00
54) 4-Nitrophenol-d4	14.742	143	19596	23.439	ng/ul	0.00
60) Fluorene-d10	15.506	176	158558	27.158	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.641	200	18420	15.529	ng/ul	0.00
73) Anthracene-d10	17.356	188	257996	29.839	ng/ul	0.00
81) Pyrene-d10	19.648	212	308864	28.945	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.383	264	332349	29.485	ng/ul	0.00
Target Compounds						
					Qvalue	
80) Fluoranthene	19.313	202	23113	1.806	ng/ul#	98
82) Pyrene	19.677	202	25140	1.892	ng/ul	98
85) Benzo(a)anthracene	21.493	228	16880	1.177	ng/ul	99
87) Chrysene	21.551	228	16278m	1.233	ng/ul	
90) Benzo(b)fluoranthene	23.625	252	19987	1.432	ng/ul	94
93) Benzo(a)pyrene	24.448	252	13405	1.011	ng/ul#	79

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

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