

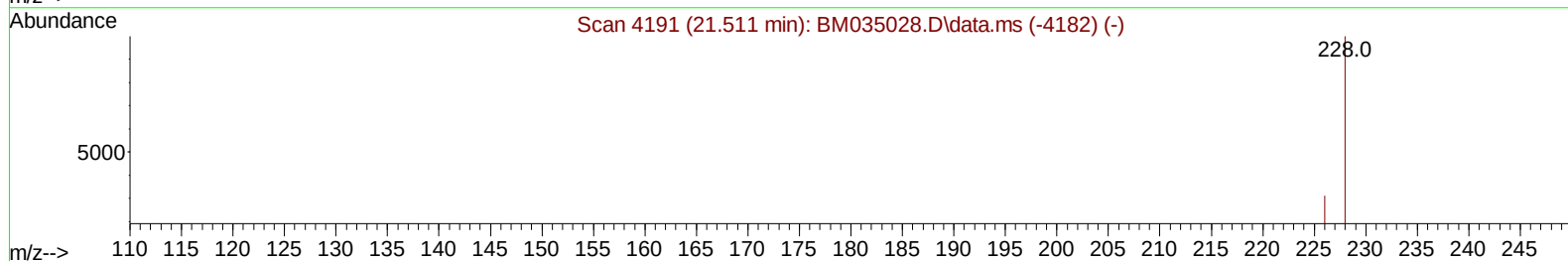
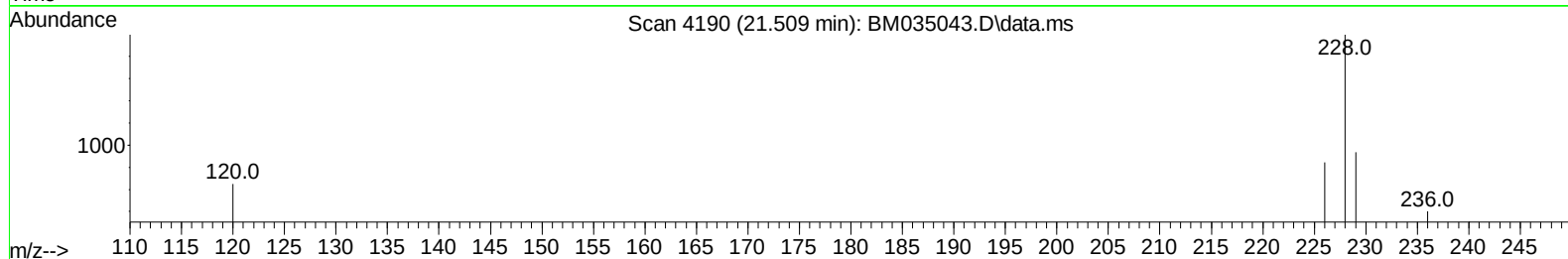
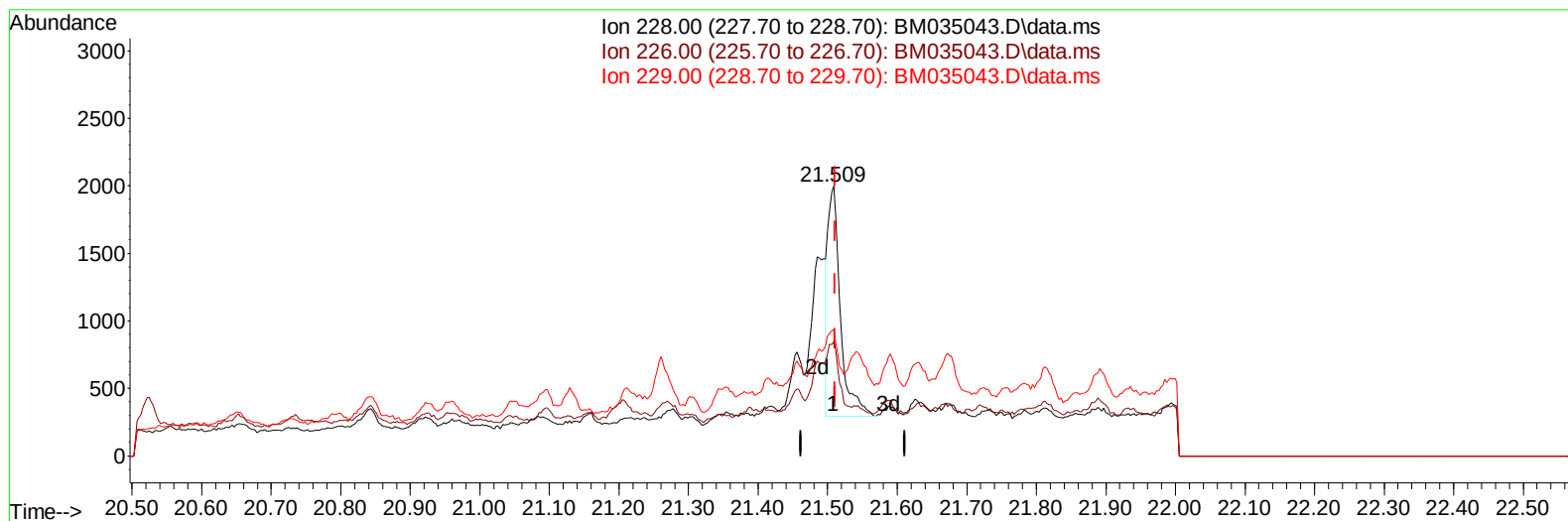
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051322\
 Data File : BM035043.D
 Acq On : 14 May 2022 00:29
 Operator : CG/JU
 Sample : N2603-18
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW498

Manual IntegrationsAPPROVED

Quant Time: May 14 01:41:46 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM051222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 12 18:15:39 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/16/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BM035043.D\data.ms

(22) Chrysene

21.509min (-0.003) 0.03 ng/u1

response 2148

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	32.30	42.36#
229.00	20.90	46.82#
0.00	0.00	0.00

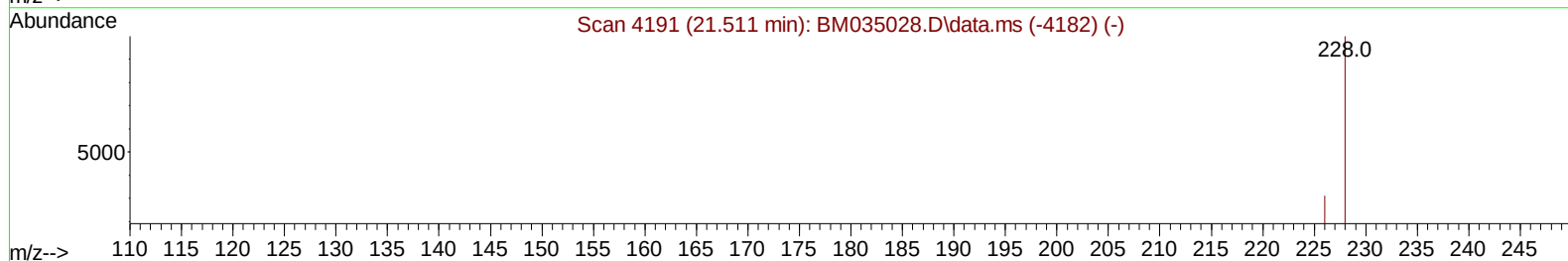
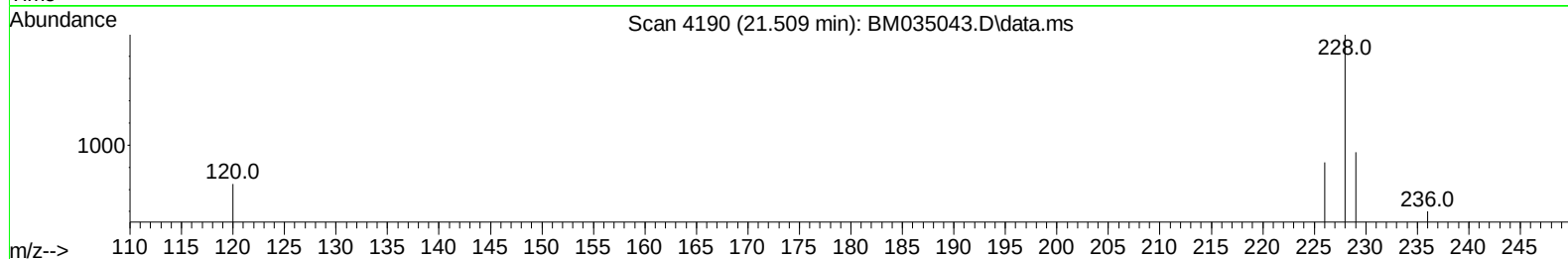
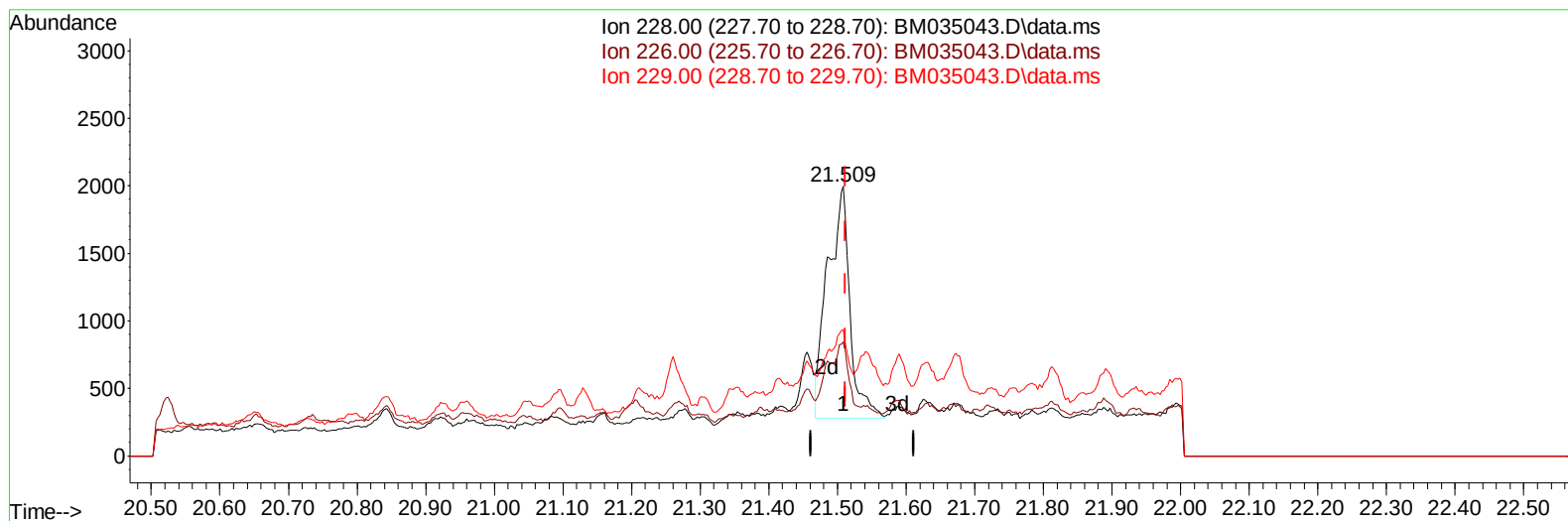
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TIC: BM035043.D\data.ms

(22) Chrysene

21.509min (-0.003) 0.05 ng/ul m

response 3880

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	32.30	42.36#
229.00	20.90	46.82#
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	7.917	152		3966	0.400	ng/ul	0.00
4) Naphthalene-d8	10.715	136		12863	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.552	164		8731	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.292	188		18732	0.400	ng/ul	0.00
17) Chrysene-d12	21.471	240		14810	0.400	ng/ul	# 0.00
23) Perylene-d12	23.856	264		14790	0.400	ng/ul	# 0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.356	96		18546	3.904	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.305	152		5788	0.282	ng/ul	0.00
18) Fluoranthene-d10	19.317	212		14794	0.291	ng/ul	0.00
Target Compounds							
						Qvalue	
5) Naphthalene	10.765	128		1371	0.032	ng/ul#	74
7) 2-Methylnaphthalene	12.382	142		3541	0.127	ng/ul	96
8) 1-Methylnaphthalene	12.596	142		2675	0.103	ng/ul	93
15) Phenanthrene	17.334	178		11966	0.168	ng/ul	98
19) Fluoranthene	19.350	202		2552	0.031	ng/ul#	6
20) Pyrene	19.712	202		3051	0.036	ng/ul#	68
22) Chrysene	21.509	228		3880m	0.054	ng/ul	
24) Benzo(b)fluoranthene	23.128	252		1724	0.022	ng/ul#	1
29) Benzo(g,h,i)perylene	27.077	276		1983	0.033	ng/ul#	58

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

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