

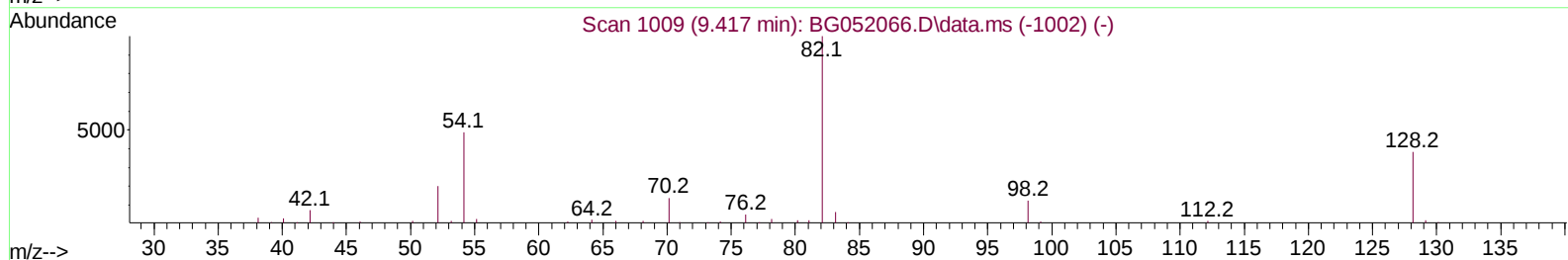
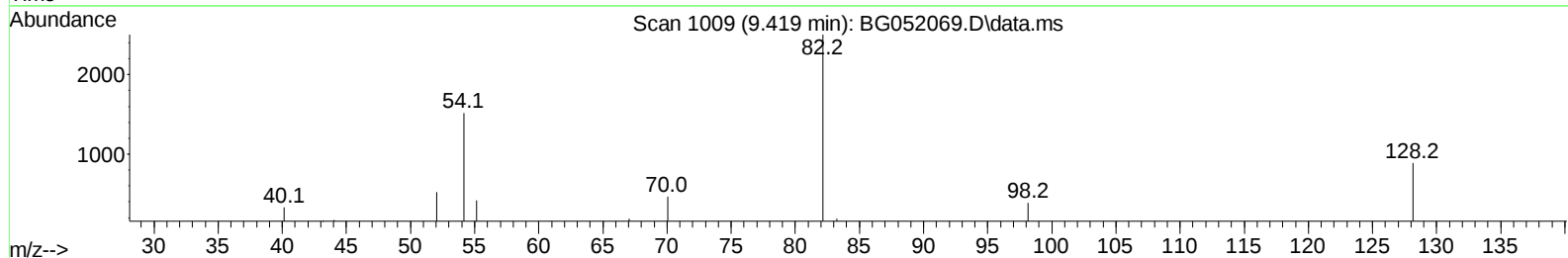
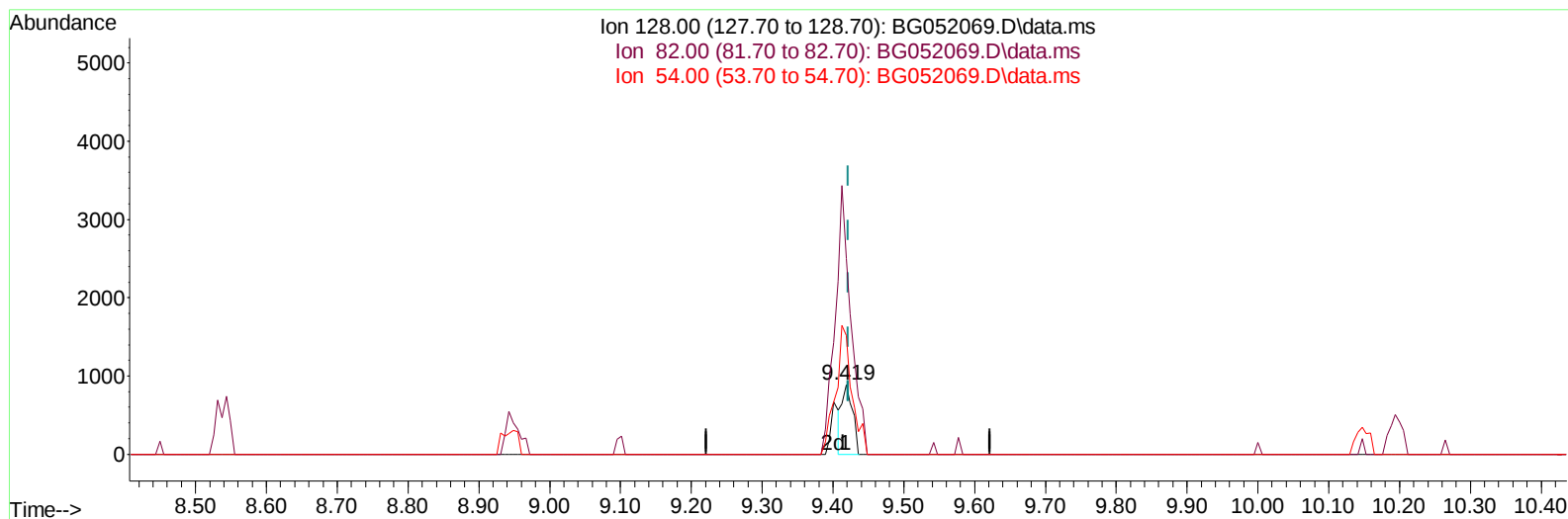
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011922\
 Data File : BG052069.D
 Acq On : 19 Jan 2022 11:37
 Operator : CG/JU
 Sample : N1100-14MEDL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLS6MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 19 23:59:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/20/2022
 Supervised By :Yogesh Patel 01/23/2022



TIC: BG052069.D\data.ms

(21) Nitrobenzene-d5 (S)

9.419min (-0.003) 0.96 ng/u1

response 942

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	265.30	281.35
54.00	131.10	170.90#
0.00	0.00	0.00

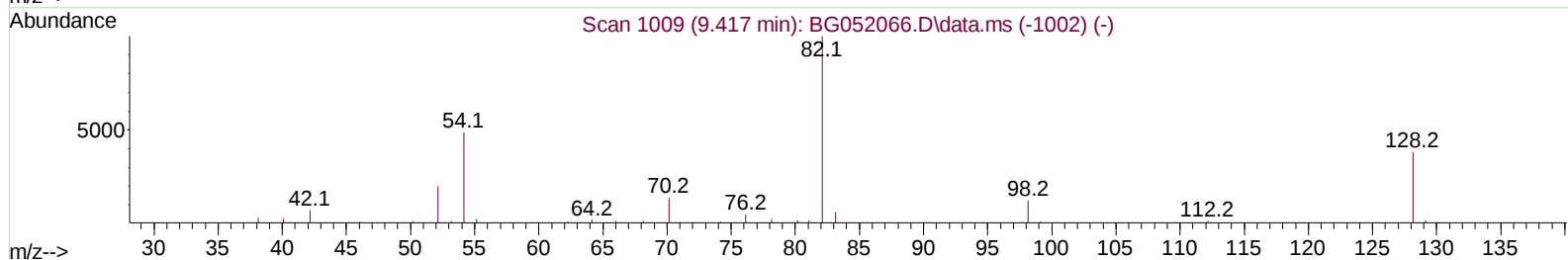
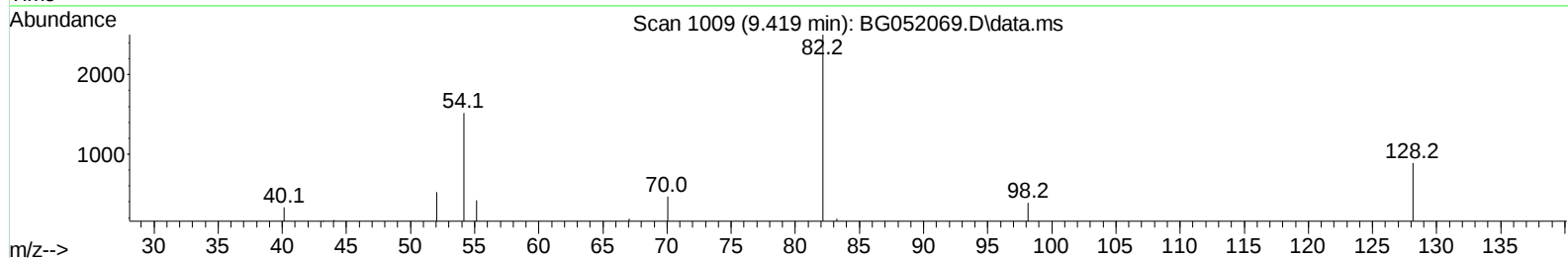
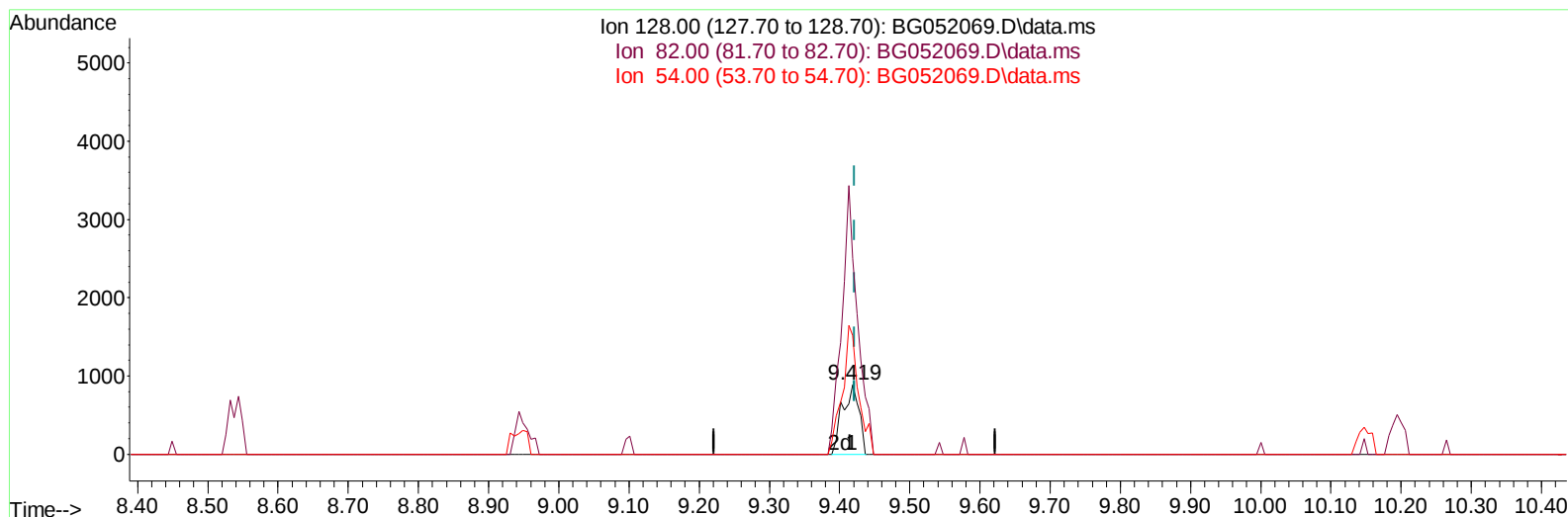
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TIC: BG052069.D\data.ms

(21) Nitrobenzene-d5 (S)

9.419min (-0.003) 1.47 ng/ul m

response 1441

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	265.30	281.35
54.00	131.10	170.90#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.244	152	27160	20.000	ng/ul	0.00
20) Naphthalene-d8	11.081	136	123743	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.888	164	93543	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.637	188	200119	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.960	240	186068	20.000	ng/ul	# 0.00
88) Perylene-d12	25.444	264	193949	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	7.392	99	4047	1.522	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.551	67	2919	1.730	ng/ul	0.00
11) 2-Chlorophenol-d4	7.774	132	2860	1.625	ng/ul	0.00
15) 4-Methylphenol-d8	8.943	113	3596	1.740	ng/ul	0.00
21) Nitrobenzene-d5	9.419	128	1441m	1.474	ng/ul	0.00
24) 2-Nitrophenol-d4	10.153	143	1732	1.602	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.693	165	3789	1.793	ng/ul	0.00
31) 4-Chloroaniline-d4	11.204	131	5279	1.793	ng/ul	0.00
46) Dimethylphthalate-d6	14.277	166	12740	1.648	ng/ul	0.00
49) Acenaphthylene-d8	14.582	160	17443	1.868	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.875	176	12540	1.860	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.737	188	20869	2.198	ng/ul	0.00
81) Pyrene-d10	20.016	212	23356	2.181	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.197	264	21370	2.093	ng/ul	0.00
Target Compounds						
					Qvalue	
86) Bis(2-ethylhexyl)phtha...	21.819	149	202162	32.601	ng/ul#	98

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

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