

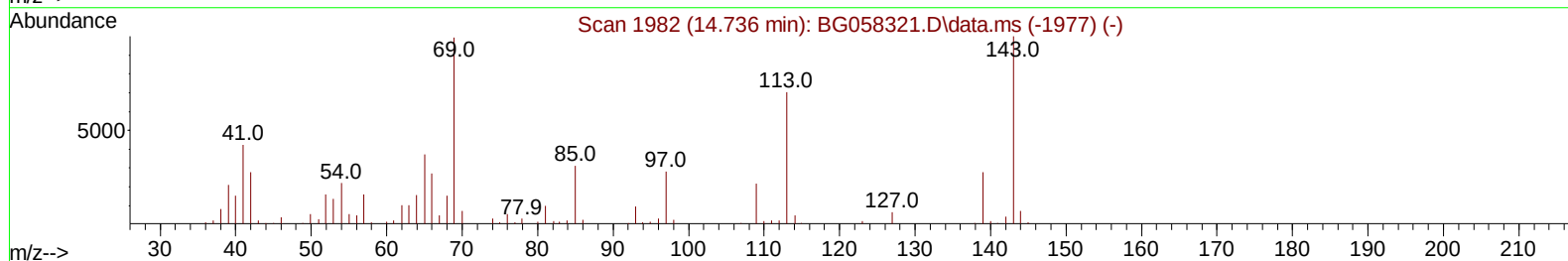
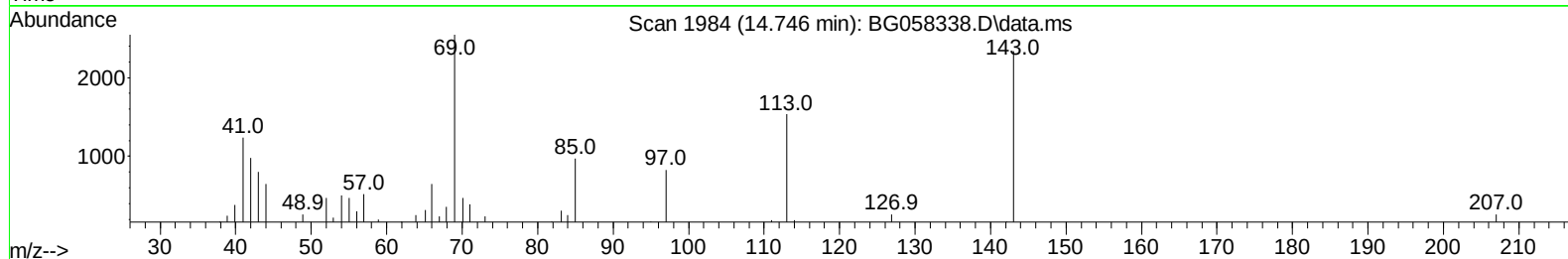
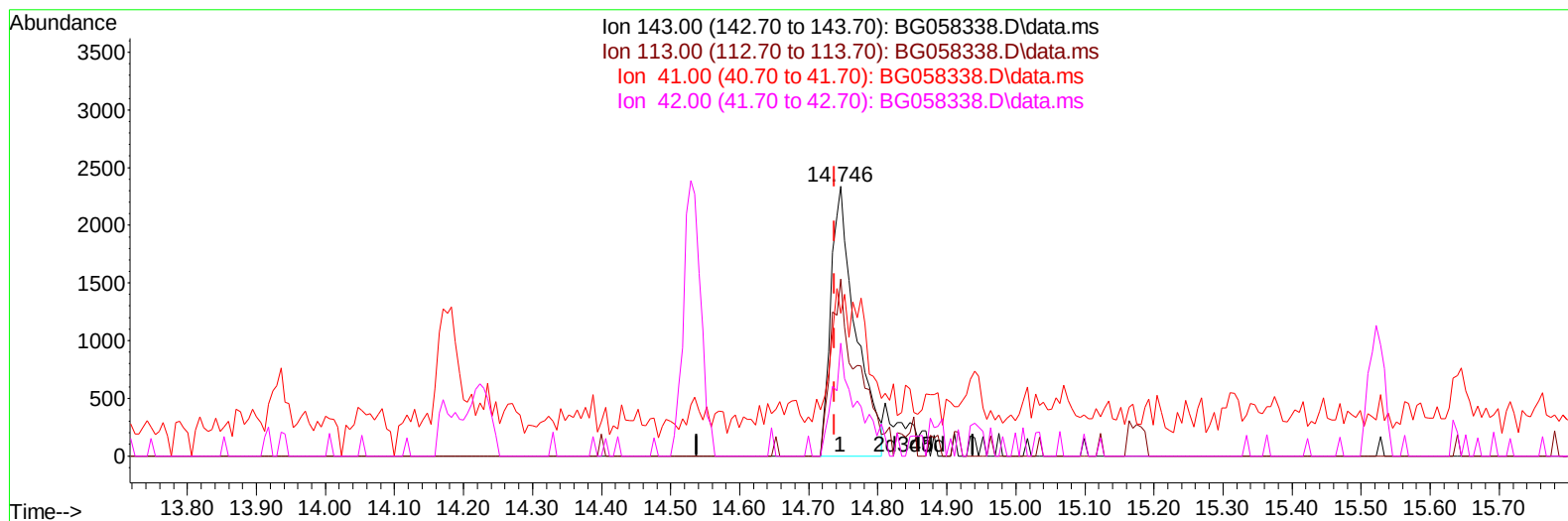
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
 Data File : BG058338.D
 Acq On : 19 Jul 2023 22:15
 Operator : CG/JU
 Sample : 03452-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A46N8

Manual IntegrationsAPPROVED

Quant Time: Jul 20 01:44:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 18 02:19:39 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/20/2023
 Supervised By :mohammad ahmed 07/20/2023



TIC: BG058338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.746min (+ 0.008) 3.87 ng/ul

response 5814

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.80	65.67
41.00	46.40	52.91
42.00	30.60	41.87#

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Quant Time: Jul 20 01:46:47 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.895	152		35825	20.000	ng/ul	0.00
20) Naphthalene-d8	10.698	136		169734	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.535	164		128985	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.279	188		303849	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240		255707	20.000	ng/ul	0.00
88) Perylene-d12	24.658	264		290695	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.336	96		1025	1.053	ng/uL	0.00
4) Pyridine-d5	3.742	84		23421	8.107	ng/ul	0.00
7) Phenol-d5	7.061	99		13456	3.959	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.226	67		16688	7.856	ng/ul	0.00
11) 2-Chlorophenol-d4	7.431	132		11711	4.923	ng/ul	0.00
15) 4-Methylphenol-d8	8.606	113		360	0.139	ng/ul	0.00
21) Nitrobenzene-d5	9.059	128		9734	7.651	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143		10735	7.729	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165		6481	2.434	ng/ul	0.00
31) 4-Chloroaniline-d4	10.839	131		213	0.055	ng/ul	0.00
46) Dimethylphthalate-d6	13.935	166		75768	7.618	ng/ul	0.00
49) Acenaphthylene-d8	14.223	160		74425	7.140	ng/ul	0.00
54) 4-Nitrophenol-d4	14.746	143		6447m	4.292	ng/ul	0.00
60) Fluorene-d10	15.522	176		70606	8.574	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200		5806	3.750	ng/ul	0.00
73) Anthracene-d10	17.373	188		36028	2.744	ng/ul	0.00
81) Pyrene-d10	19.664	212		100379	6.646	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.441	264		15151	1.078	ng/ul	0.00
Target Compounds							
						Qvalue	
16) Acetophenone	8.718	105		11726	2.656	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.409	149		19484	1.984	ng/ul	97

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

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