

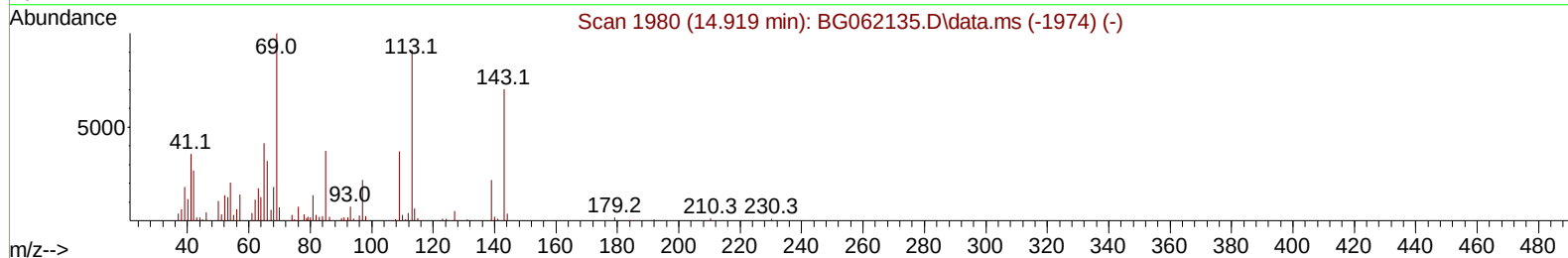
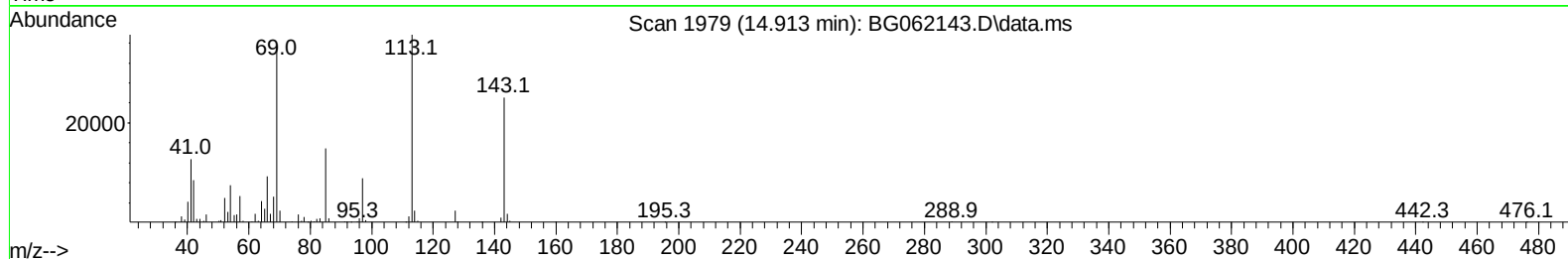
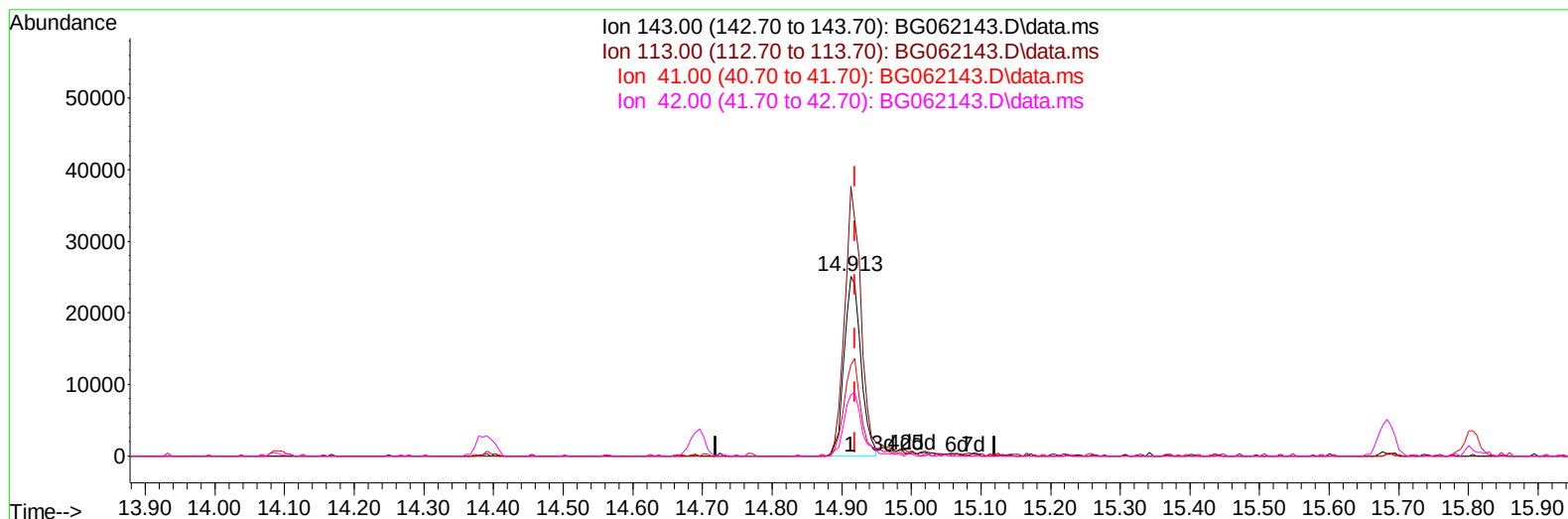
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071924\
Data File : BG062143.D
Acq On : 19 Jul 2024 16:54
Operator : MA/JU
Sample : PB162080BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK080

Manual IntegrationsAPPROVED

Quant Time: Jul 19 18:59:52 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 19 14:38:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/20/2024
Supervised By :mohammad ahmed 07/23/2024



TIC: BG062143.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.913min (-0.006) 27.16 ng/ul

response 42140

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	129.40	149.98
41.00	50.80	50.90
42.00	38.10	34.20

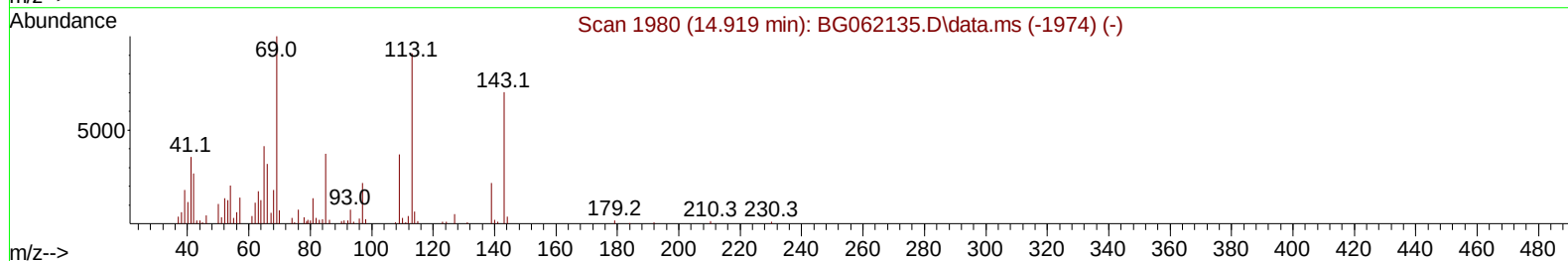
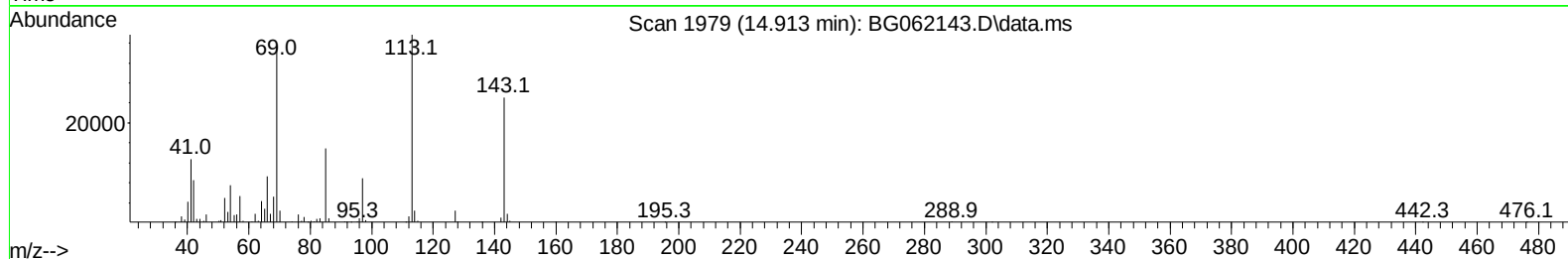
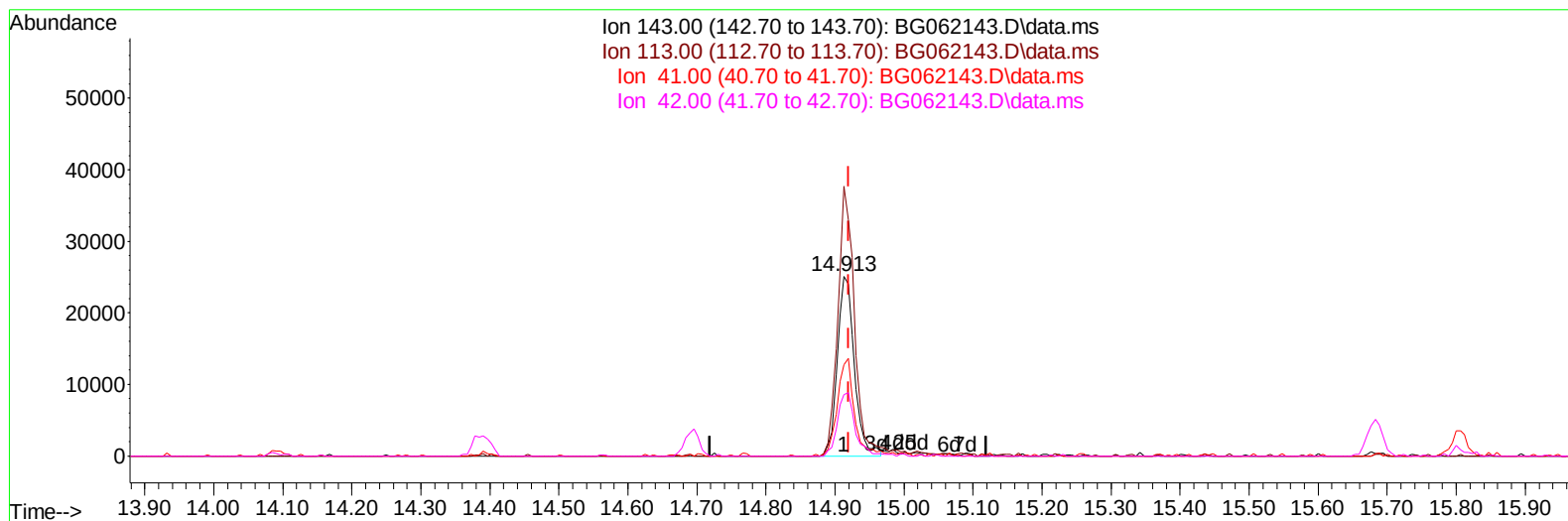
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TIC: BG062143.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.913min (-0.006) 27.82 ng/ul m

response 43155

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	129.40	149.98
41.00	50.80	50.90
42.00	38.10	34.20

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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 07/23/2024

Quant Time: Jul 20 02:12:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 19 14:38:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.064	152	44174	20.000	ng/ul	0.00
20) Naphthalene-d8	10.877	136	175450	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.696	164	151891	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	435795	20.000	ng/ul	0.00
79) Chrysene-d12	21.698	240	495768	20.000	ng/ul	0.00
88) Perylene-d12	24.929	264	501650	20.000	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.458	96	6731	7.114	ng/uL	0.00
4) Pyridine-d5	3.875	84	89704	26.094	ng/ul	0.00
7) Phenol-d5	7.235	99	110627	25.438	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.388	67	68997	27.613	ng/ul	0.00
11) 2-Chlorophenol-d4	7.600	132	69658	25.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.780	113	84240	25.434	ng/ul	0.00
21) Nitrobenzene-d5	9.233	128	38098	27.255	ng/ul	0.00
24) 2-Nitrophenol-d4	9.955	143	46865	27.855	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.501	165	96286	27.338	ng/ul	0.00
31) 4-Chloroaniline-d4	11.018	131	110333	26.042	ng/ul	0.00
46) Dimethylphthalate-d6	14.091	166	390895	30.362	ng/ul	0.00
49) Acenaphthylene-d8	14.390	160	373924	28.644	ng/ul	0.00
54) 4-Nitrophenol-d4	14.913	143	43155m	27.816	ng/ul	0.00
60) Fluorene-d10	15.683	176	319688	29.164	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.806	200	78555	28.365	ng/ul	0.00
73) Anthracene-d10	17.539	188	595280	29.242	ng/ul	0.00
81) Pyrene-d10	19.818	212	832088	29.602	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.706	264	764313	30.479	ng/ul	-0.01

Target Compounds Qvalue

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

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