

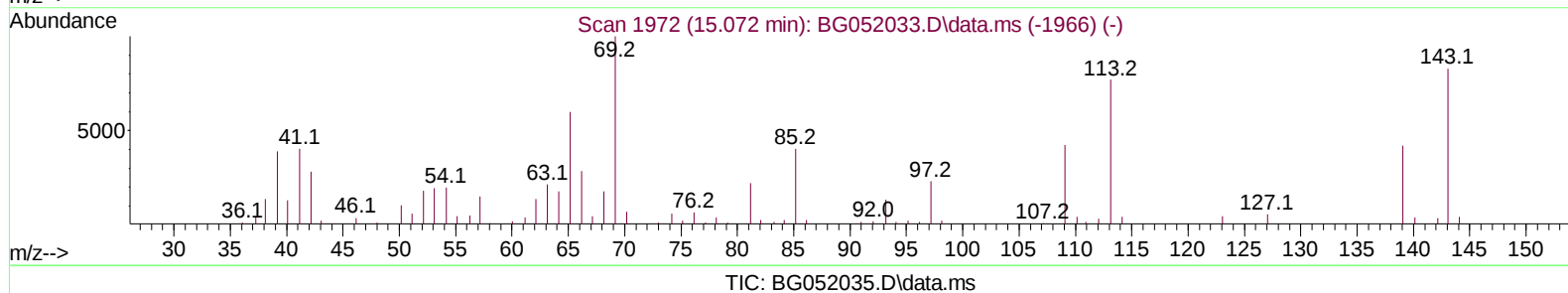
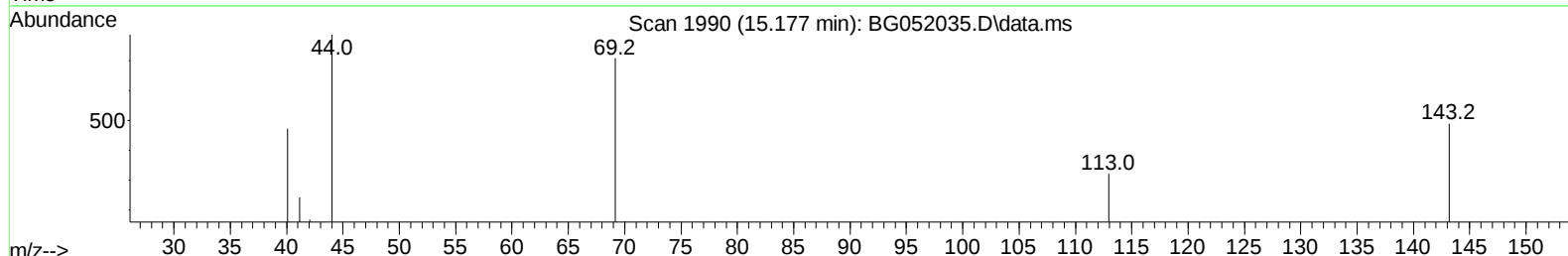
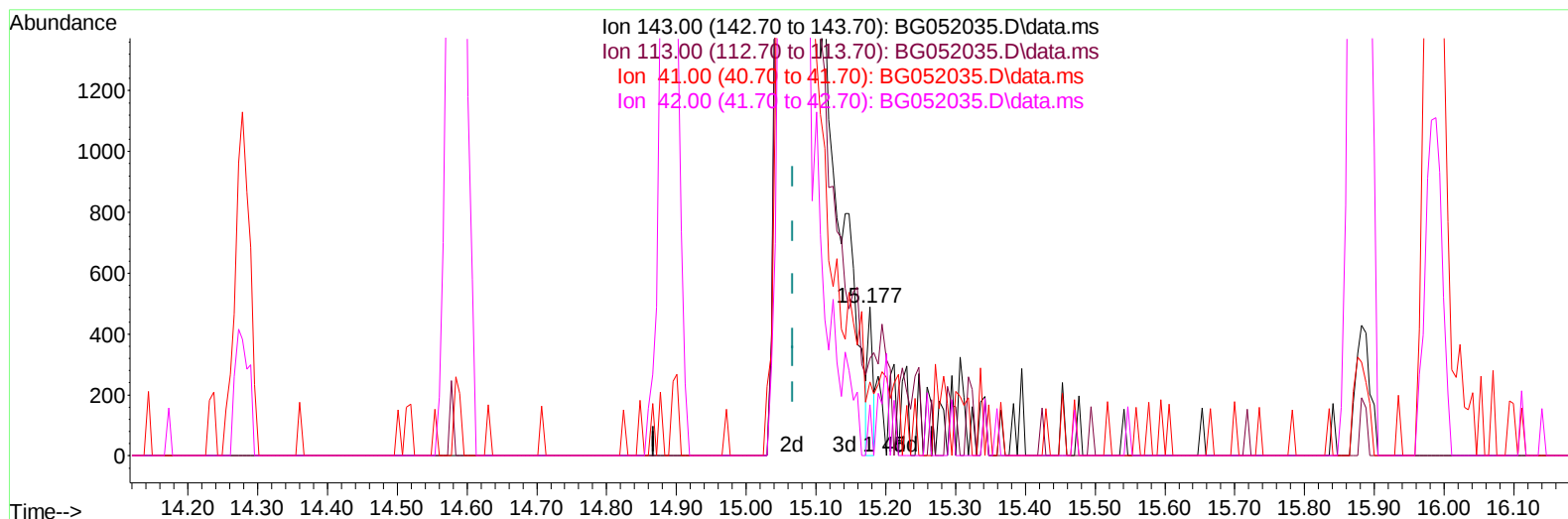
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052035.D
 Acq On : 17 Jan 2022 22:32
 Operator : CG/JU
 Sample : PB142057BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK057

Manual IntegrationsAPPROVED

Quant Time: Jan 18 00:03:25 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 01/18/2022
 Supervised By :mohammad ahmed 01/21/2022



(54) 4-Nitrophenol-d4 (S)

15.177min (+ 0.110) 0.13 ng/ul

response 246

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	65.64
41.00	44.40	49.69
42.00	29.70	34.36

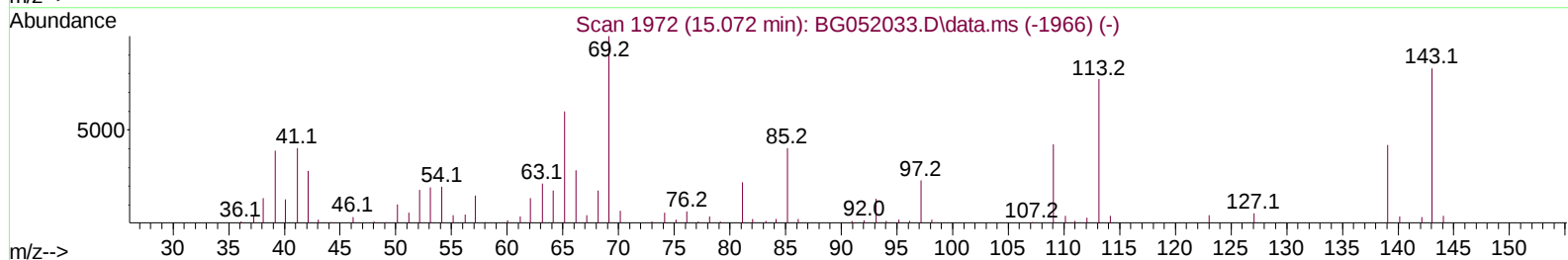
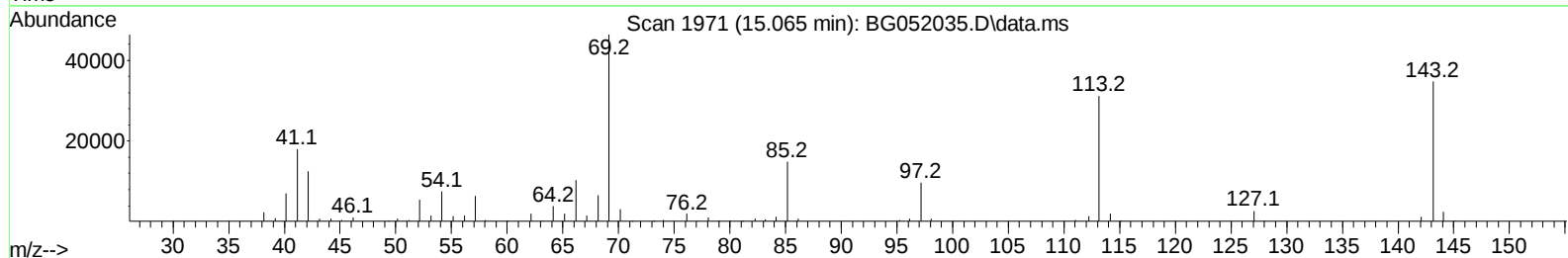
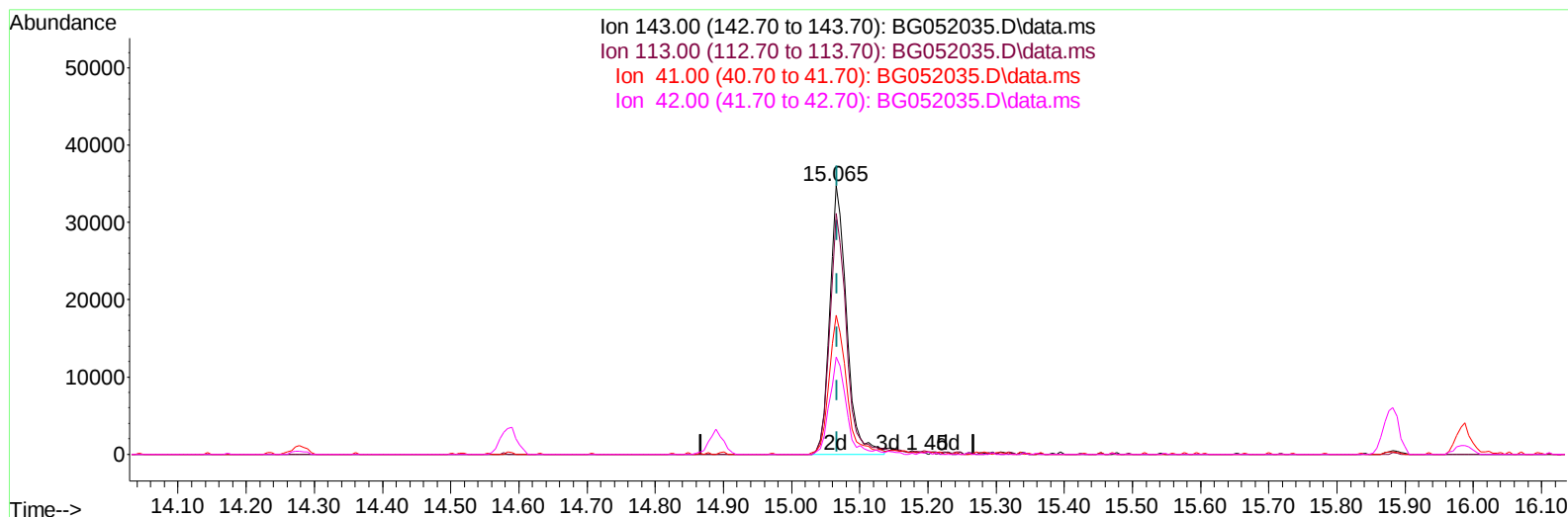
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TIC: BG052035.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.065min (-0.002) 31.96 ng/ul m

response 59550

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	89.80
41.00	44.40	51.74
42.00	29.70	36.15#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.245	152	40063	20.000	ng/ul	0.00
20) Naphthalene-d8	11.088	136	184165	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.889	164	134847	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.638	188	316619	20.000	ng/ul	0.00
79) Chrysene-d12	21.962	240	279751	20.000	ng/ul	0.00
88) Perylene-d12	25.463	264	270574	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.575	96	7531	6.717	ng/uL	0.00
4) Pyridine-d5	3.998	84	110035	31.571	ng/ul	0.00
7) Phenol-d5	7.382	99	131740	33.598	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.552	67	87526	35.163	ng/ul	0.00
11) 2-Chlorophenol-d4	7.769	132	85817	33.056	ng/ul	0.00
15) 4-Methylphenol-d8	8.944	113	102639	33.673	ng/ul	0.00
21) Nitrobenzene-d5	9.420	128	47333	32.526	ng/ul	0.00
24) 2-Nitrophenol-d4	10.149	143	53444	33.207	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.695	165	93518	29.727	ng/ul	0.00
31) 4-Chloroaniline-d4	11.212	131	136872	31.242	ng/ul	0.00
46) Dimethylphthalate-d6	14.278	166	351994	31.580	ng/ul	0.00
49) Acenaphthylene-d8	14.584	160	427959	31.786	ng/ul	0.00
54) 4-Nitrophenol-d4	15.065	143	59550m	31.955	ng/ul	0.00
60) Fluorene-d10	15.882	176	302569	31.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.988	200	54095	27.083	ng/ul	0.00
73) Anthracene-d10	17.738	188	525956	35.018	ng/ul	0.00
81) Pyrene-d10	20.018	212	596440	37.046	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.216	264	511709	35.924	ng/ul	0.02

Target Compounds Qvalue

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

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