

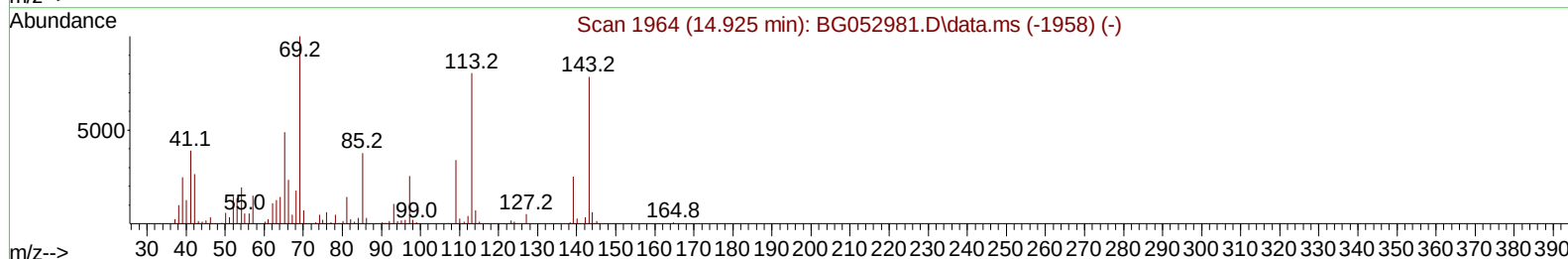
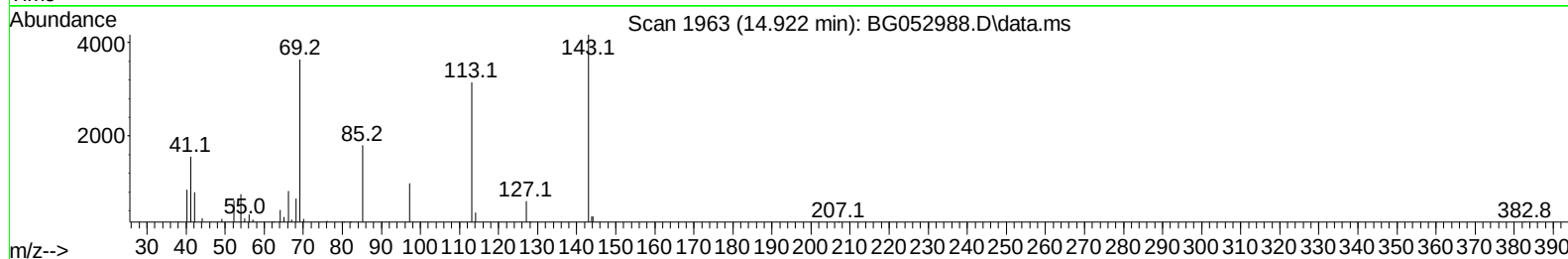
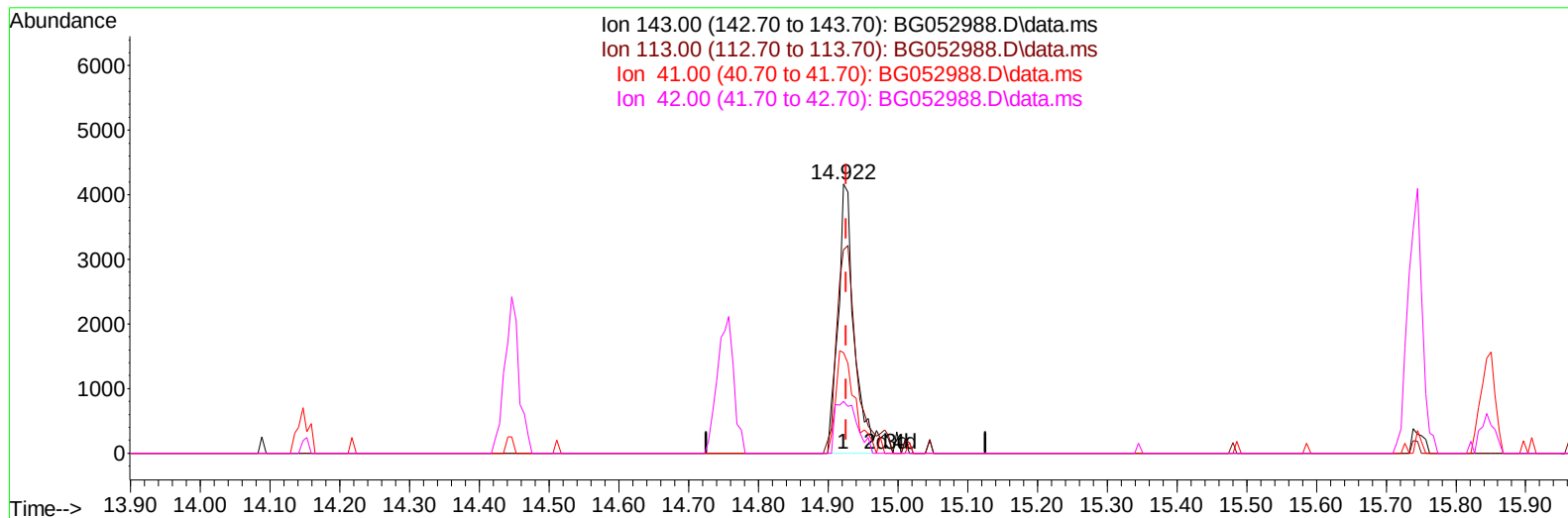
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052988.D
Acq On : 31 Mar 2022 20:35
Operator : CG/JU
Sample : N2120-01RX
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0HS8RX

Manual IntegrationsAPPROVED

Quant Time: Apr 11 07:11:00 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/01/2022
Supervised By :Yogesh Patel 04/06/2022



TIC: BG052988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (-0.004) 5.29 ng/u1

response 6547

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	75.67#
41.00	47.20	37.32#
42.00	29.40	19.31#

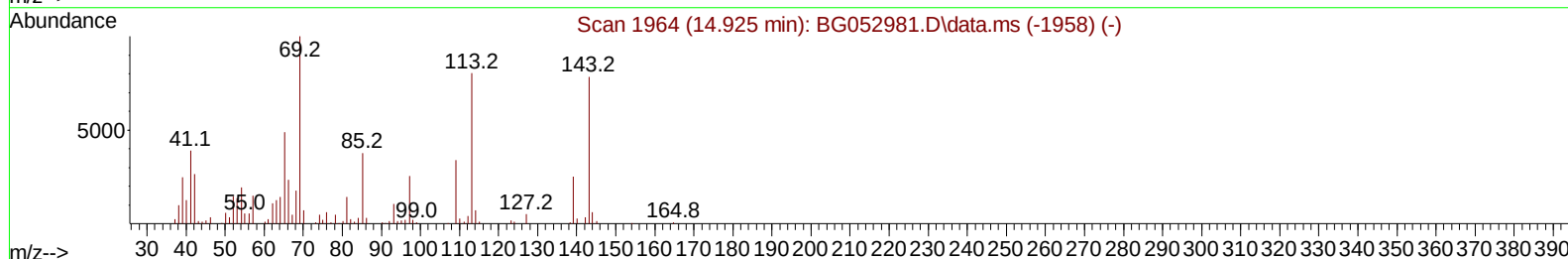
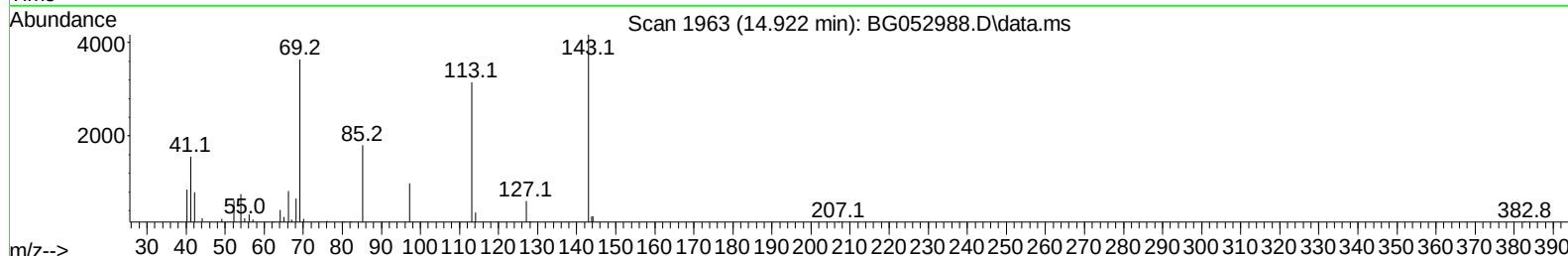
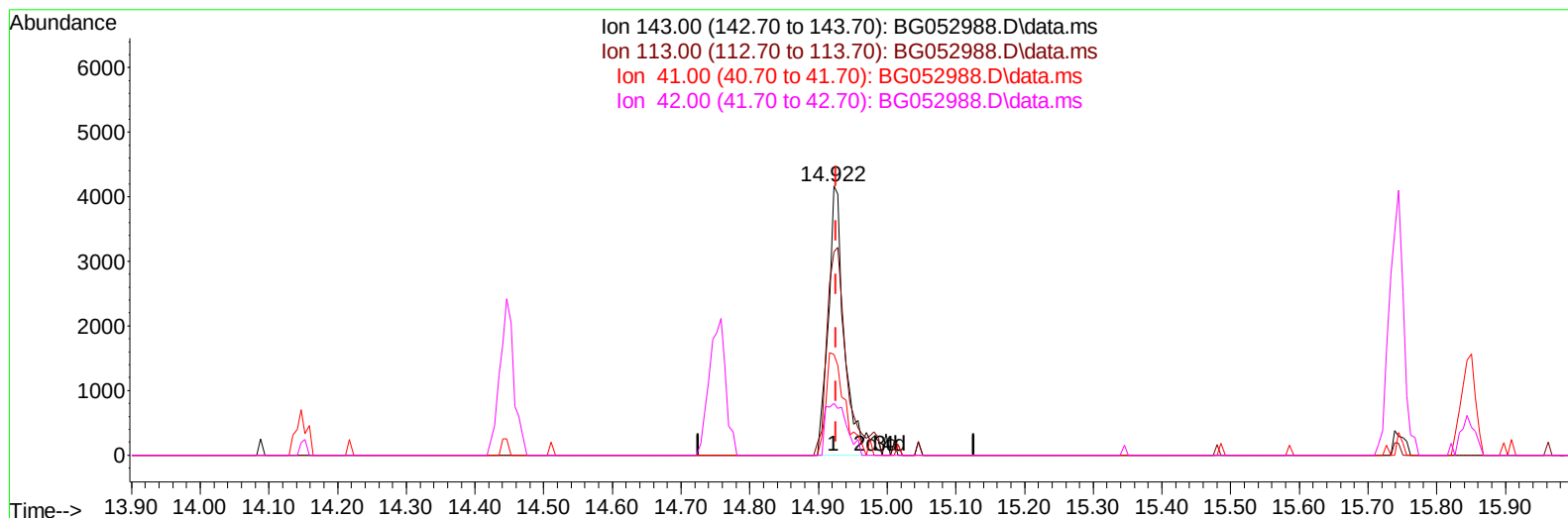
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 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0HS8RX

Manual IntegrationsAPPROVED

Quant Time: Apr 01 01:05:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 31 00:50:49 2022
 Response via : Initial Calibration

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 Supervised By :Yogesh Patel 04/06/2022



TIC: BG052988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (-0.004) 5.60 ng/ul m

response 6932

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	75.67#
41.00	47.20	37.32#
42.00	29.40	19.31#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.137	152	26835	20.000	ng/ul	0.00
20) Naphthalene-d8	10.945	136	118874	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.752	164	95803	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.495	188	230641	20.000	ng/ul	0.00
79) Chrysene-d12	21.778	240	250672	20.000	ng/ul	0.00
88) Perylene-d12	25.085	264	247580	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.549	96	3322	4.373	ng/uL	0.00
4) Pyridine-d5	3.978	84	9841	5.189	ng/ul	0.01
7) Phenol-d5	7.280	99	15646	6.510	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.450	67	47639	32.453	ng/ul	0.00
11) 2-Chlorophenol-d4	7.667	132	39106	23.962	ng/ul	0.00
15) 4-Methylphenol-d8	8.825	113	27707	15.914	ng/ul	0.00
21) Nitrobenzene-d5	9.294	128	28050	32.780	ng/ul	0.00
24) 2-Nitrophenol-d4	10.017	143	29067	32.205	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	52208	26.589	ng/ul	0.00
31) 4-Chloroaniline-d4	11.069	131	45922	17.696	ng/ul	0.00
46) Dimethylphthalate-d6	14.153	166	224363	30.484	ng/ul	0.00
49) Acenaphthylene-d8	14.446	160	248518	27.788	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	6932m	5.602	ng/ul	0.00
60) Fluorene-d10	15.745	176	201819	31.229	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.844	200	31493	26.832	ng/ul	0.00
73) Anthracene-d10	17.595	188	347703	32.374	ng/ul	0.00
81) Pyrene-d10	19.874	212	427803	30.507	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.856	264	416086	32.526	ng/ul	0.00

Target Compounds Qvalue

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

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