

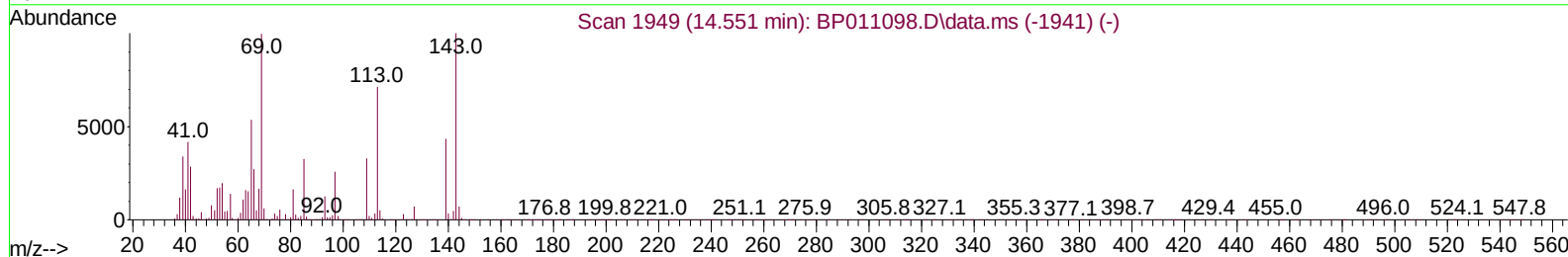
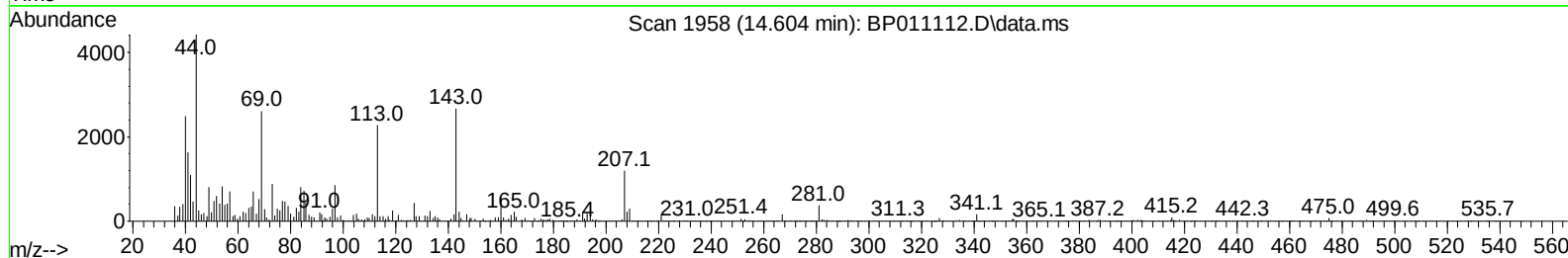
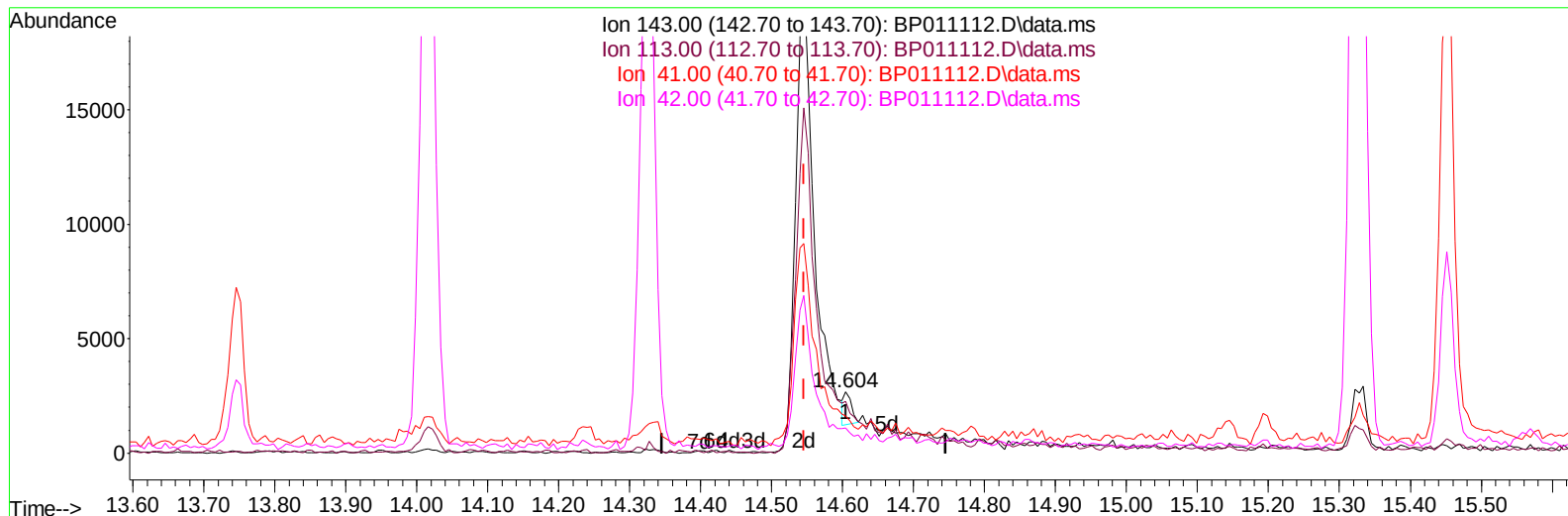
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
Data File : BP011112.D
Acq On : 26 Jul 2022 16:13
Operator : CG/JU
Sample : N3809-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YBE68

Manual IntegrationsAPPROVED

Quant Time: Jul 26 16:56:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 26 14:31:36 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/27/2022
Supervised By :mohammad ahmed 07/27/2022



TIC: BP011112.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.604min (+ 0.059) 0.07 ng/ul

response 1049

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	85.14
41.00	59.00	61.14
42.00	41.10	40.96

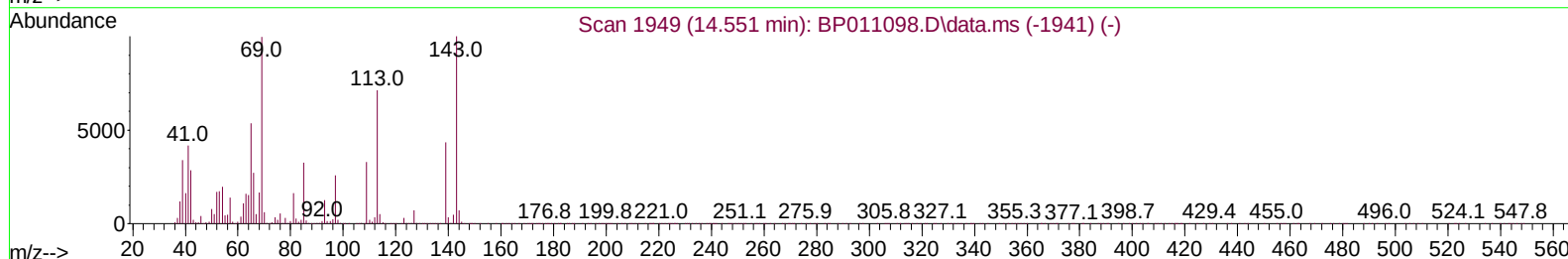
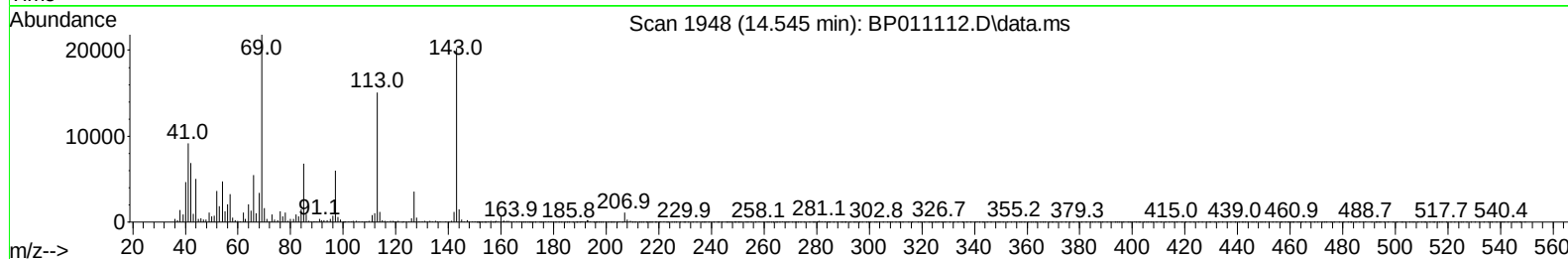
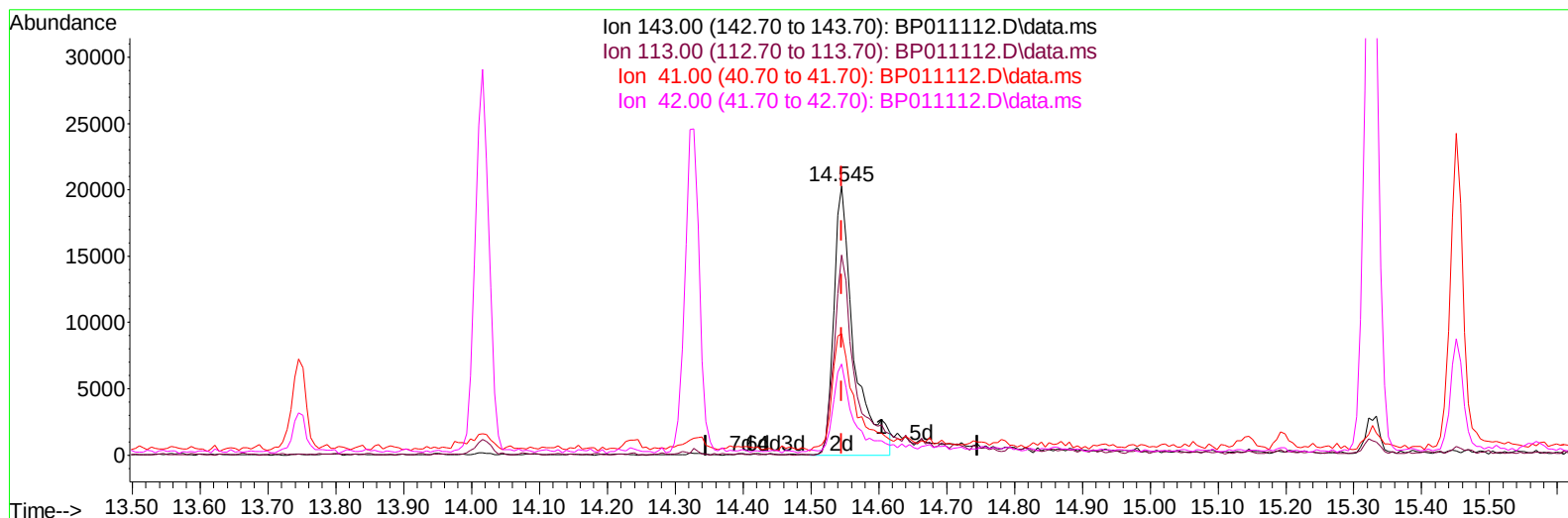
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TIC: BP011112.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.000) 2.96 ng/ul m

response 42221

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	74.40#
41.00	59.00	45.18#
42.00	41.10	33.96

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	418413	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	1864315	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.328	164	1031576	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1761030	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	1299580	20.000	ng/ul	# 0.00
88) Perylene-d12	23.557	264	1379076	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	61261	5.278	ng/uL	0.00
4) Pyridine-d5	3.575	84	285810	9.527	ng/ul	0.00
7) Phenol-d5	6.840	99	246340	7.172	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	785011	35.544	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132	730310	26.315	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	471070	17.132	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	481550	35.444	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	468755	34.868	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	734780	29.887	ng/ul	0.00
31) 4-Chloroaniline-d4	10.604	131	1055326	26.037	ng/ul	0.00
46) Dimethylphthalate-d6	13.746	166	2653707	37.639	ng/ul	0.00
49) Acenaphthylene-d8	14.016	160	3088842	37.131	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	42221m	2.959	ng/ul	0.00
60) Fluorene-d10	15.328	176	2144952	35.791	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	275854	31.656	ng/ul	0.00
73) Anthracene-d10	17.192	188	2950099	39.405	ng/ul	0.00
81) Pyrene-d10	19.451	212	2927029	42.884	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	2650939	40.023	ng/ul	0.00

Target Compounds Qvalue

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

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