

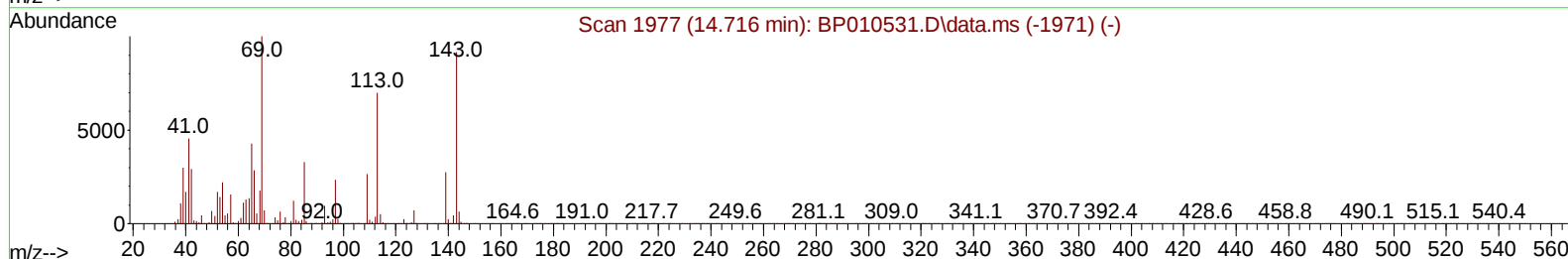
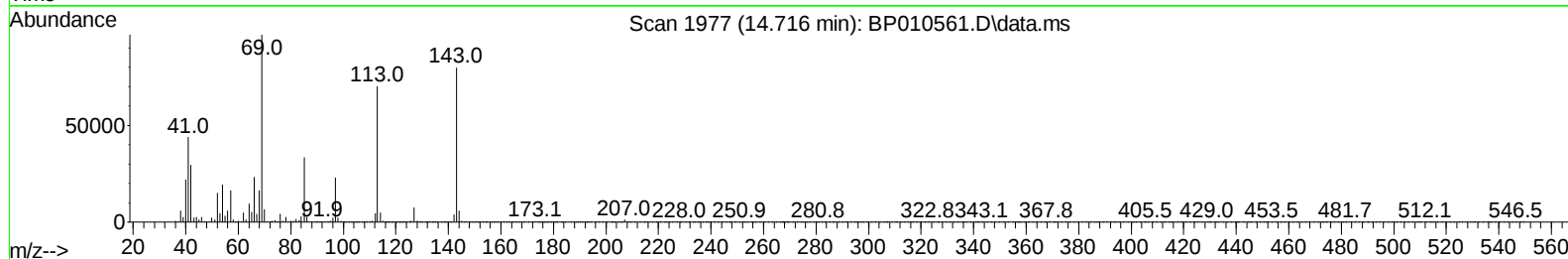
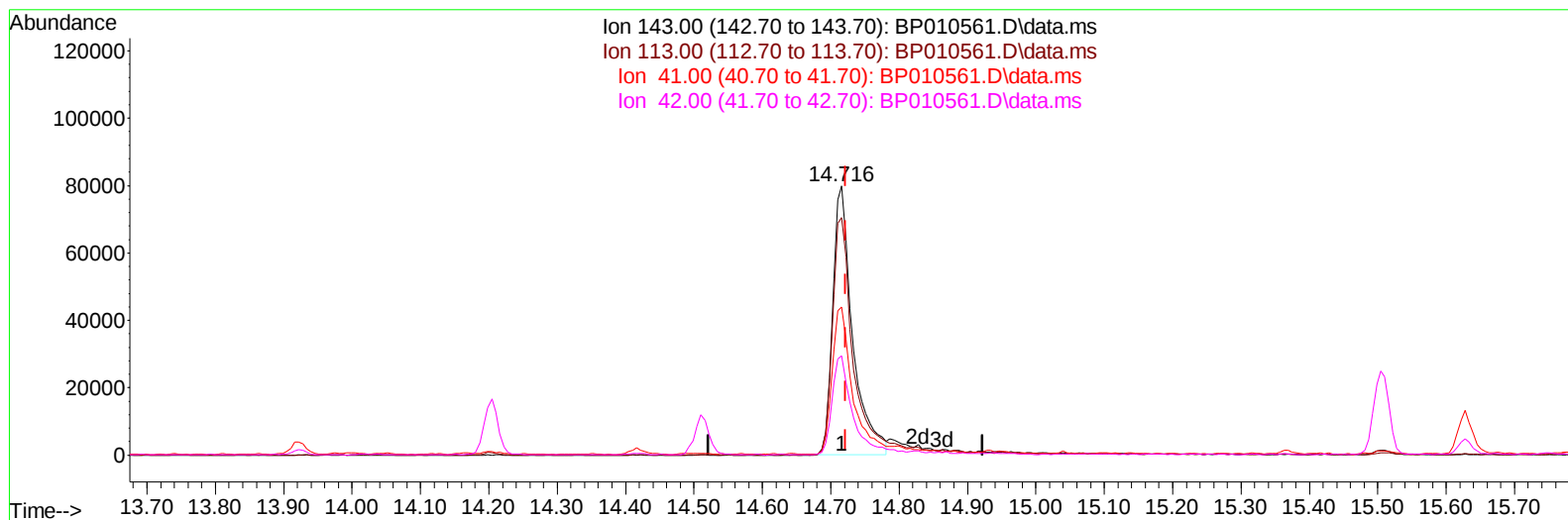
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010561.D
Acq On : 26 May 2022 05:58
Operator : CG/JU
Sample : N2883-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4K8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/26/2022
Supervised By :mohammad ahmed 05/27/2022

Quant Time: May 26 07:04:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 24 05:50:40 2022
Response via : Initial Calibration



TIC: BP010561.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.716min (-0.006) 30.59 ng/u1****response 161844**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	88.19#
41.00	23.20	55.08#
42.00	18.00	36.91#

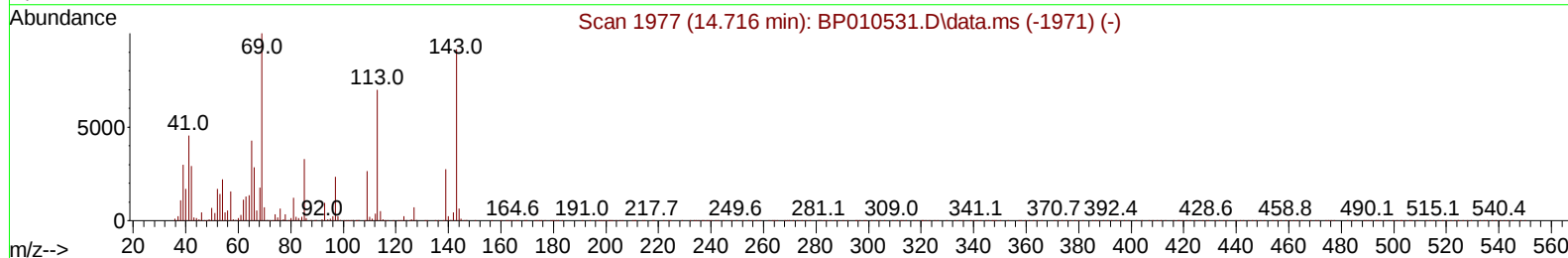
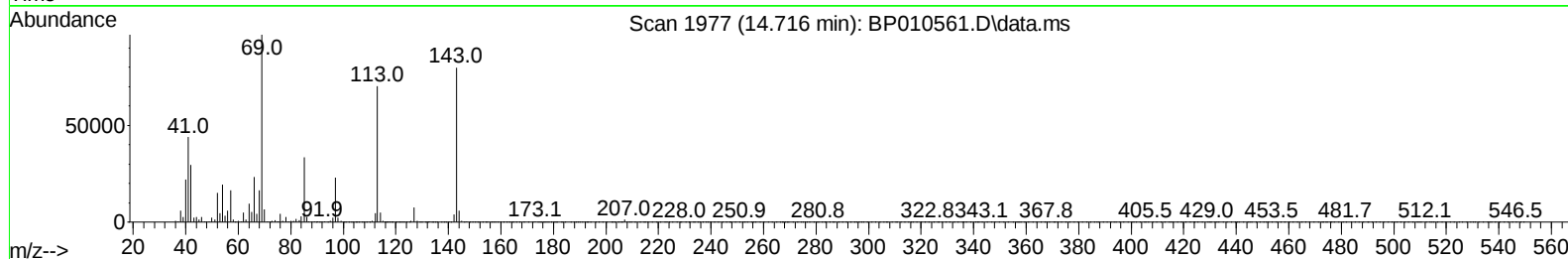
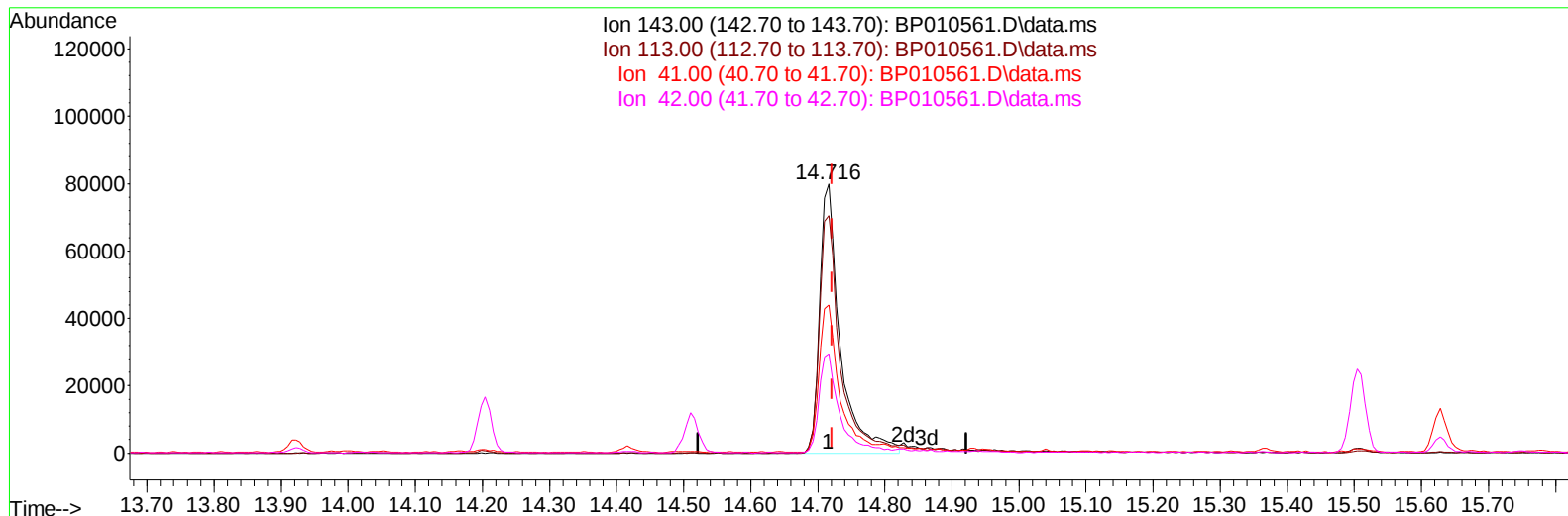
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TIC: BP010561.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (-0.006) 32.21 ng/ul m

response 170459

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	88.19#
41.00	23.20	55.08#
42.00	18.00	36.91#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	133297	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	573352	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.510	164	403496	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.263	188	903806	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.345	240	879099	20.000	ng/ul	# 0.00
88) Perylene-d12	23.763	264	774340	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	20450	6.539	ng/uL	0.00
4) Pyridine-d5	3.752	84	301978	34.009	ng/ul	0.00
7) Phenol-d5	7.046	99	402434	35.970	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	294156	39.133	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	327030	38.200	ng/ul	-0.01
15) 4-Methylphenol-d8	8.587	113	359862	37.848	ng/ul	-0.01
21) Nitrobenzene-d5	9.040	128	184973	41.685	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	201513	42.852	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	340142	38.253	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	376584	28.484	ng/ul	-0.01
46) Dimethylphthalate-d6	13.922	166	1345784	45.269	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	1505267	43.897	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	170459m	32.214	ng/ul	0.00
60) Fluorene-d10	15.504	176	1142638	44.180	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	162584	30.759	ng/ul	-0.01
73) Anthracene-d10	17.363	188	1822617	44.535	ng/ul	0.00
81) Pyrene-d10	19.598	212	2237883	47.935	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1783679	45.874	ng/ul	0.00

Target Compounds Qvalue

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

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