

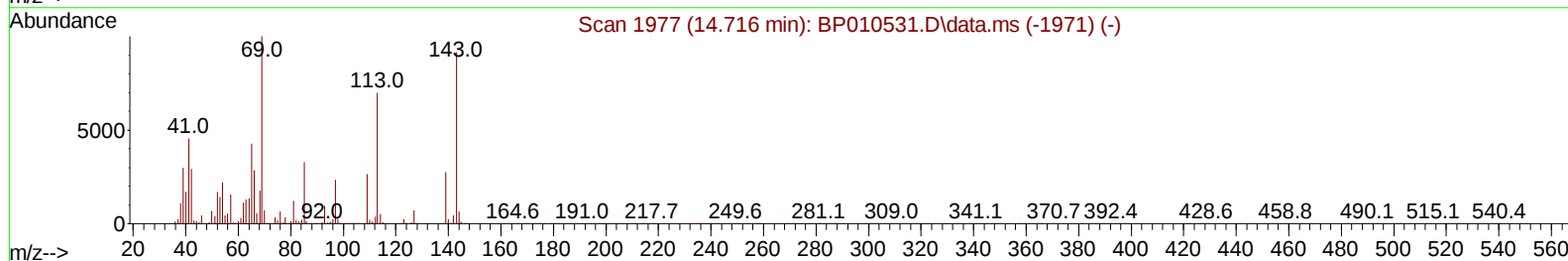
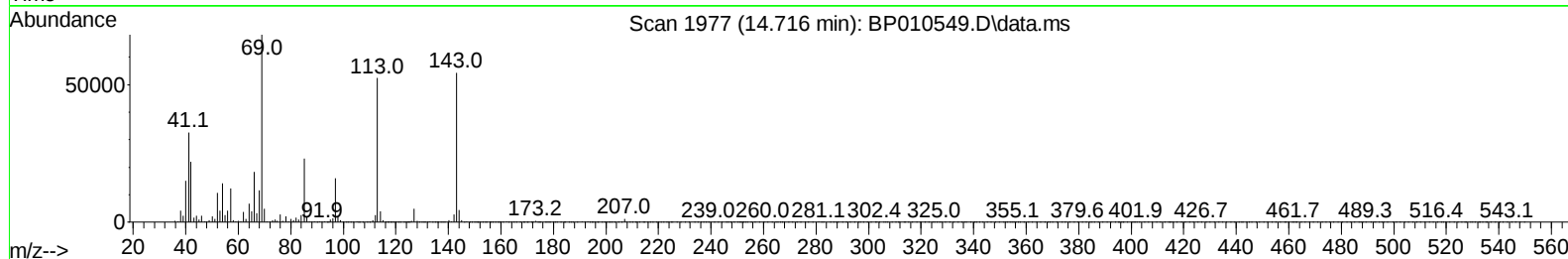
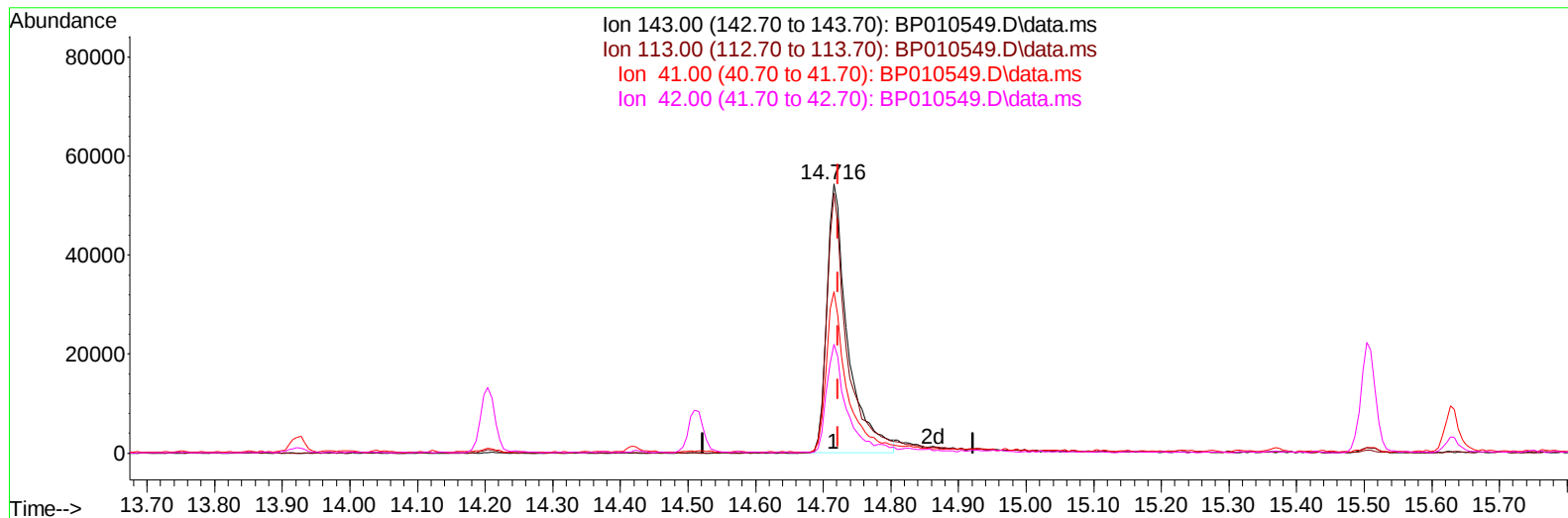
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
 Data File : BP010549.D
 Acq On : 25 May 2022 22:39
 Operator : CG/JU
 Sample : N2883-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4K3

Manual IntegrationsAPPROVED

Quant Time: May 25 23:28:47 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

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



TIC: BP010549.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (-0.006) 29.30 ng/u1

response 117678

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	96.51#
41.00	23.20	60.13#
42.00	18.00	40.37#

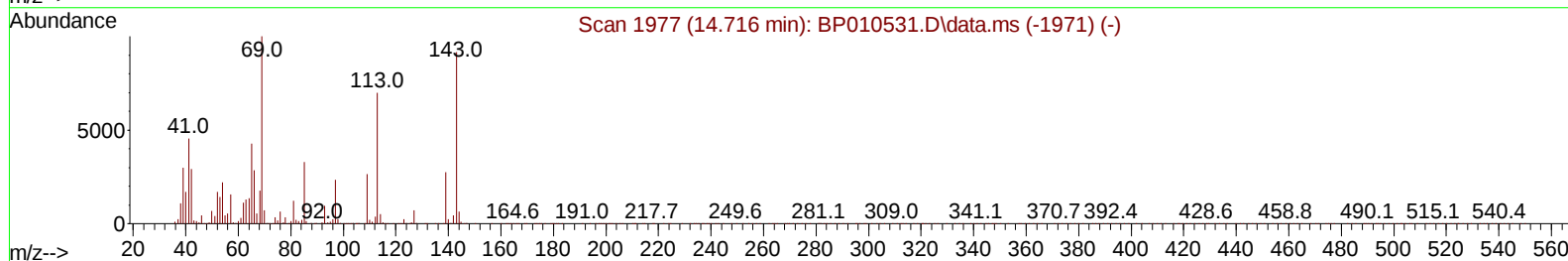
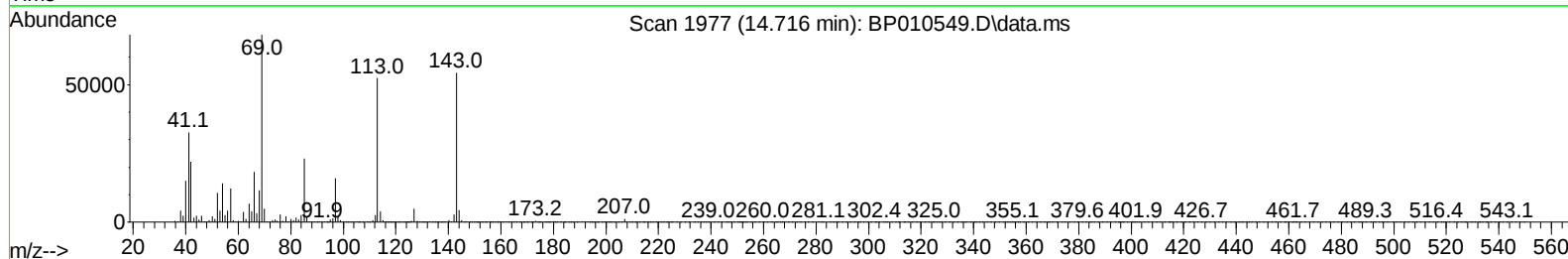
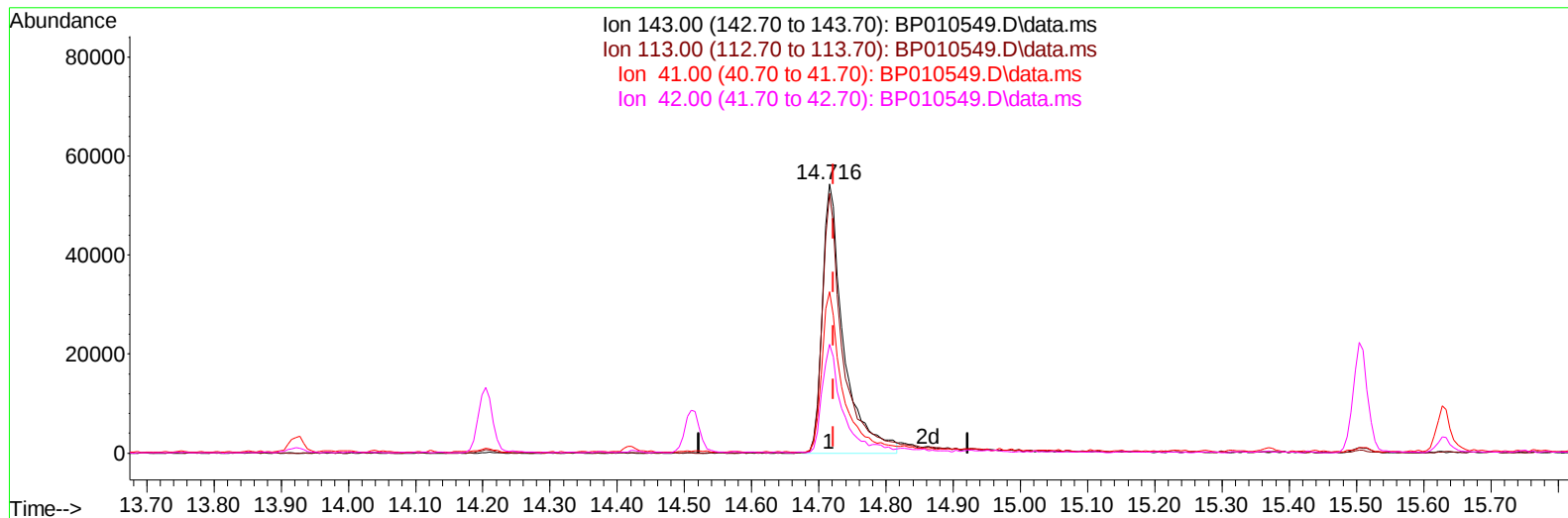
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TIC: BP010549.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (-0.006) 29.75 ng/ul m

response 119489

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	96.51#
41.00	23.20	60.13#
42.00	18.00	40.37#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	100843	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	433138	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.510	164	306283	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.263	188	703208	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.345	240	732740	20.000	ng/ul	# 0.00
88) Perylene-d12	23.762	264	673770	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	19205	8.118	ng/uL	0.00
4) Pyridine-d5	3.752	84	225990	33.642	ng/ul	0.00
7) Phenol-d5	7.046	99	313575	37.048	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	243721	42.858	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	263627	40.705	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	274699	38.190	ng/ul	0.00
21) Nitrobenzene-d5	9.040	128	149808	44.689	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	156676	44.103	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	264801	39.421	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	291501	29.186	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	1059775	46.963	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	1189230	45.688	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	119489m	29.749	ng/ul	0.00
60) Fluorene-d10	15.510	176	923581	47.045	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	117601	28.595	ng/ul	0.00
73) Anthracene-d10	17.369	188	1475717	46.344	ng/ul	0.00
81) Pyrene-d10	19.598	212	1850388	47.551	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.609	264	1627488	48.104	ng/ul	0.00

Target Compounds Qvalue

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

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