

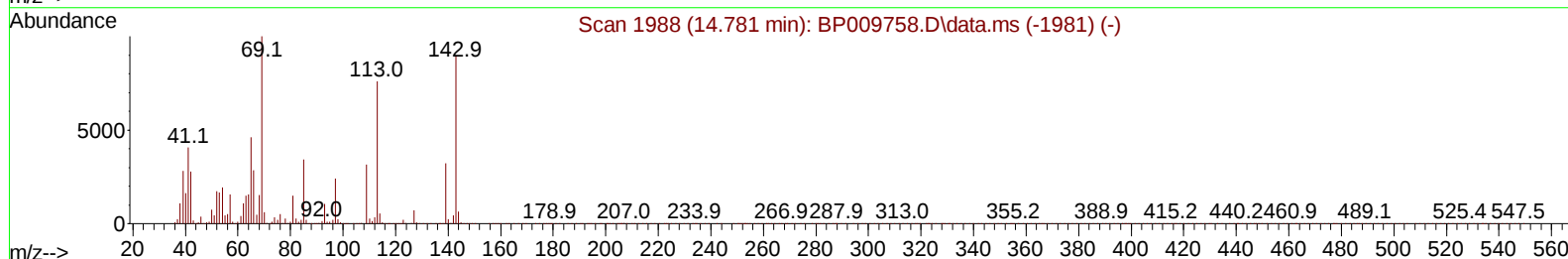
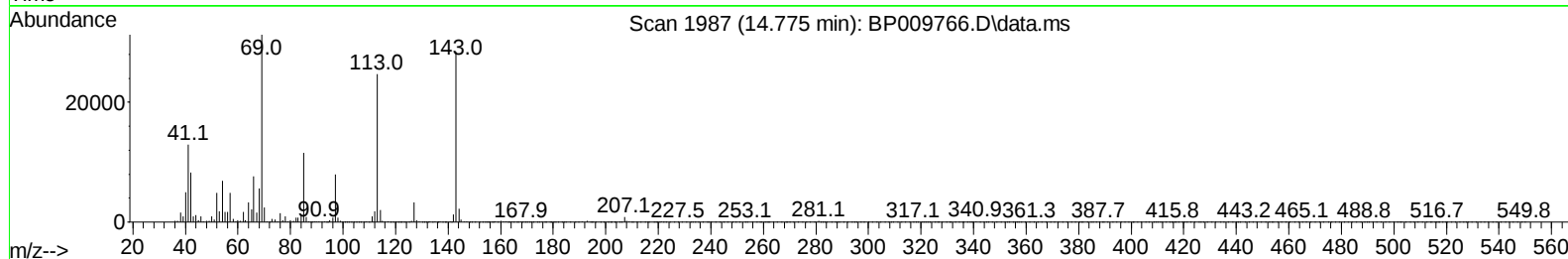
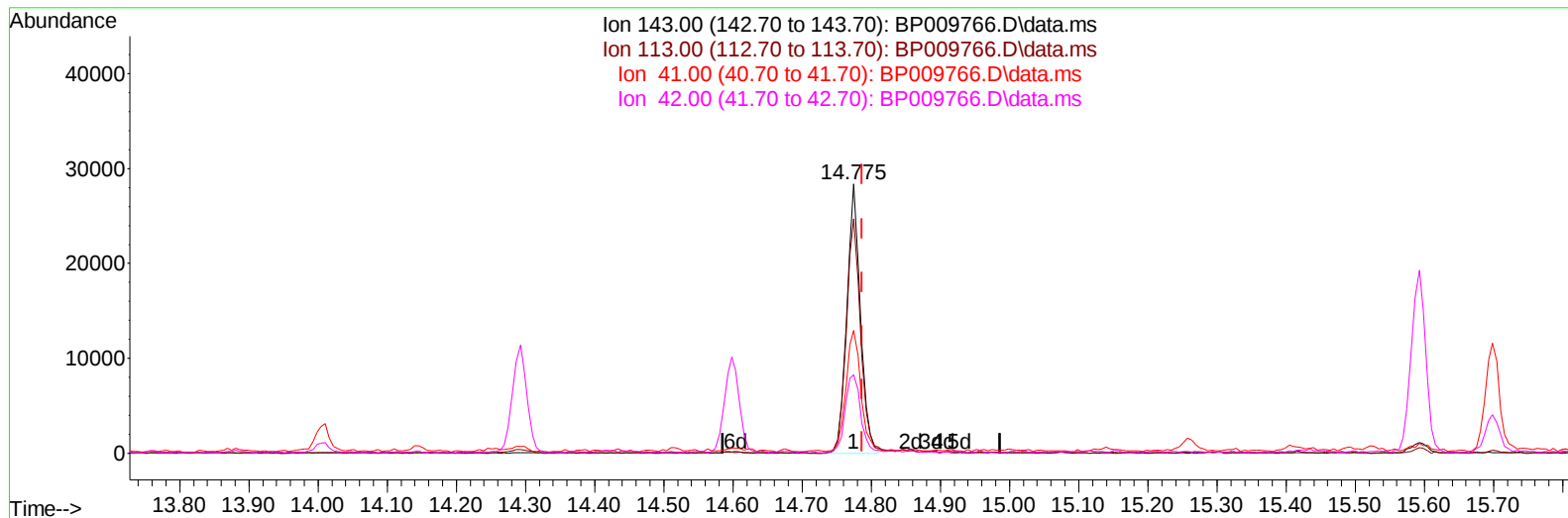
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009766.D
 Acq On : 11 Apr 2022 23:49
 Operator : CG/JU
 Sample : N2289-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFC6

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:54:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009766.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 6.63 ng/u1

response 38752

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.14#
41.00	23.20	45.54#
42.00	18.00	29.04#

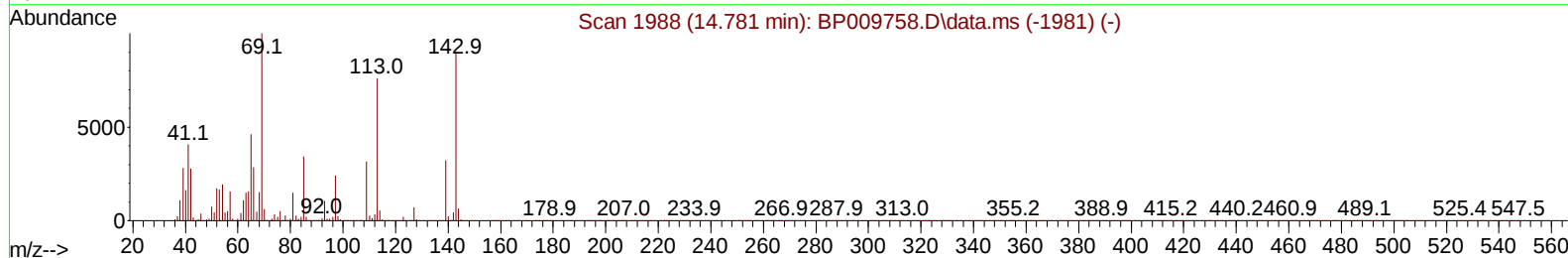
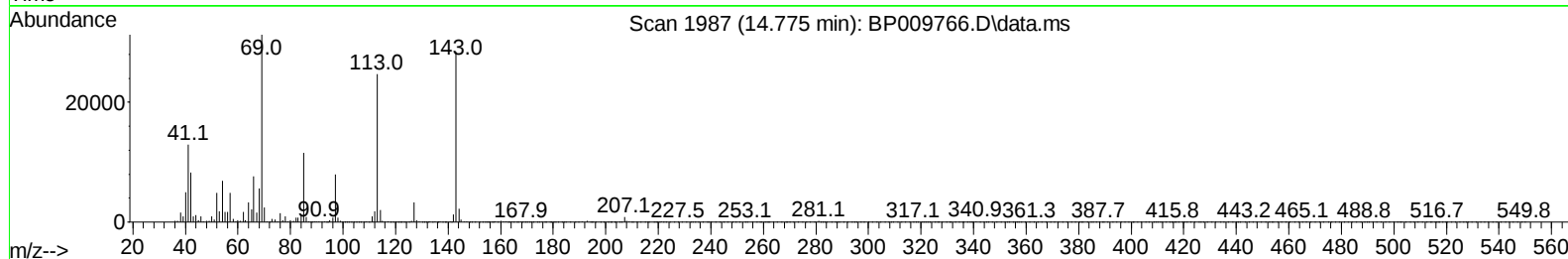
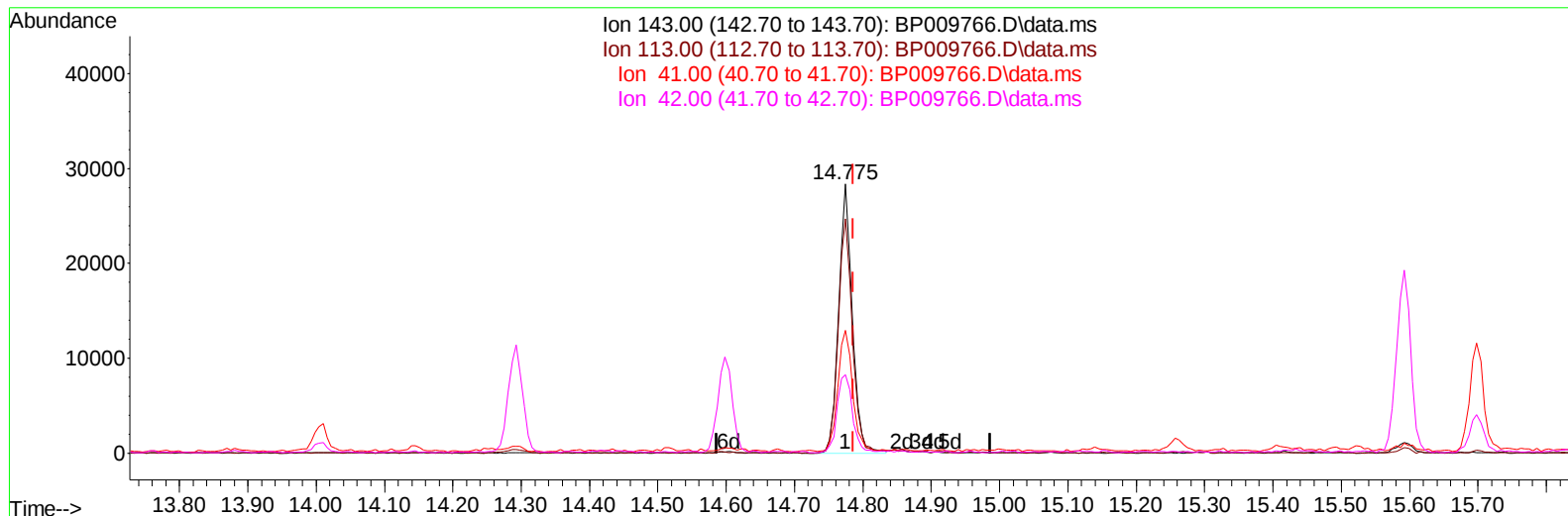
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TIC: BP009766.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 6.71 ng/ul m

response 39168

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.14#
41.00	23.20	45.54#
42.00	18.00	29.04#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	136576	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	590884	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.598	164	409990	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	918326	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.427	240	958934	20.000	ng/ul	#-0.02
88) Perylene-d12	23.898	264	897369	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	18385	6.008	ng/uL	0.00
4) Pyridine-d5	3.805	84	73985	8.769	ng/ul	0.00
7) Phenol-d5	7.122	99	75733	6.316	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	230106	32.235	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	219834	24.306	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	154420	15.427	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	145627	34.176	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	151134	32.352	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	266560	27.268	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.904	131	360025	25.858	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	1114670	34.611	ng/ul	-0.01
49) Acenaphthylene-d8	14.293	160	1182802	31.001	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	39168m	6.706	ng/ul	-0.01
60) Fluorene-d10	15.592	176	946349	34.931	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	171704	31.562	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1621032	36.868	ng/ul	-0.01
81) Pyrene-d10	19.675	212	1984546	37.893	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.739	264	1786240	38.223	ng/ul	-0.02

Target Compounds Qvalue

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

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