

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011961.D
 Acq On : 06 Oct 2022 09:55
 Operator : CG/JU
 Sample : N4956-07
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

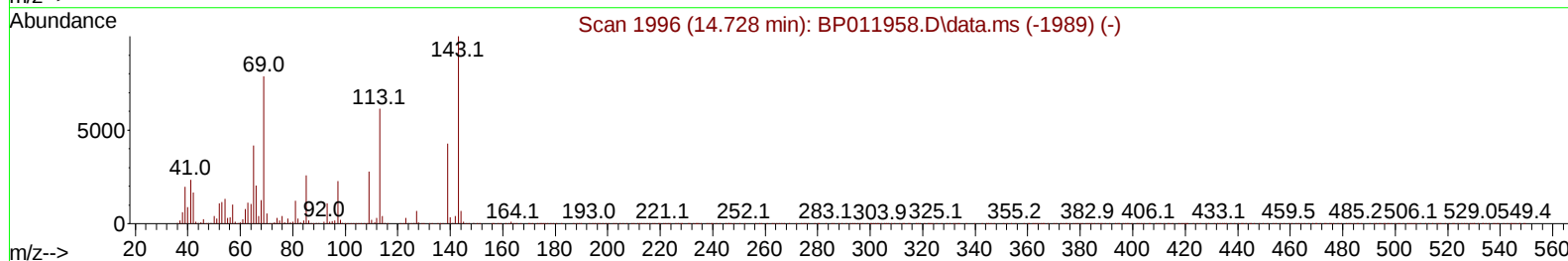
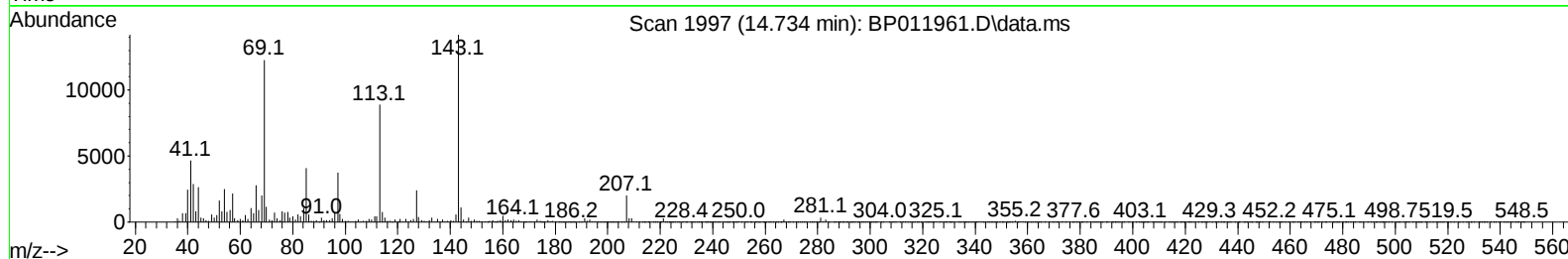
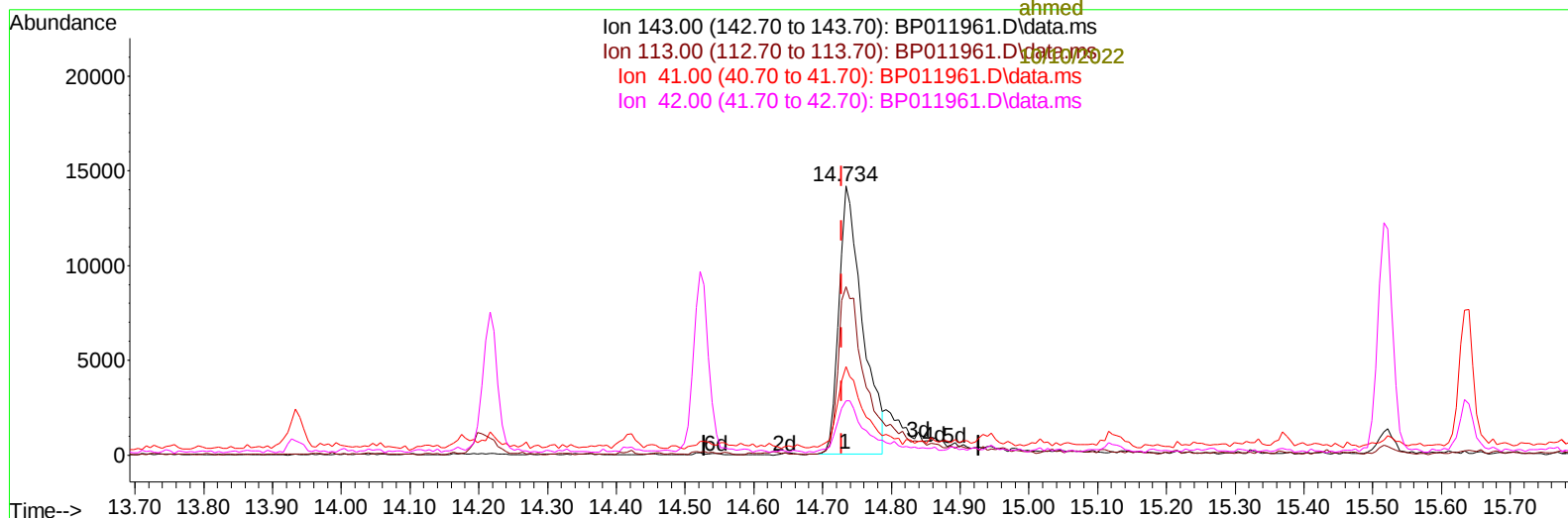
Instrument :
 BNA_P
 ClientSampleId :
 CB8S8

Manual IntegrationsAPPROVED

Quant Time: Oct 06 23:04:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

10/07/2022
 Supervised By :mohammad
 ahmed



TIC: BP011961.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 2.93 ng/u1

response 33295

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	62.70
41.00	24.20	32.80#
42.00	16.40	20.21#

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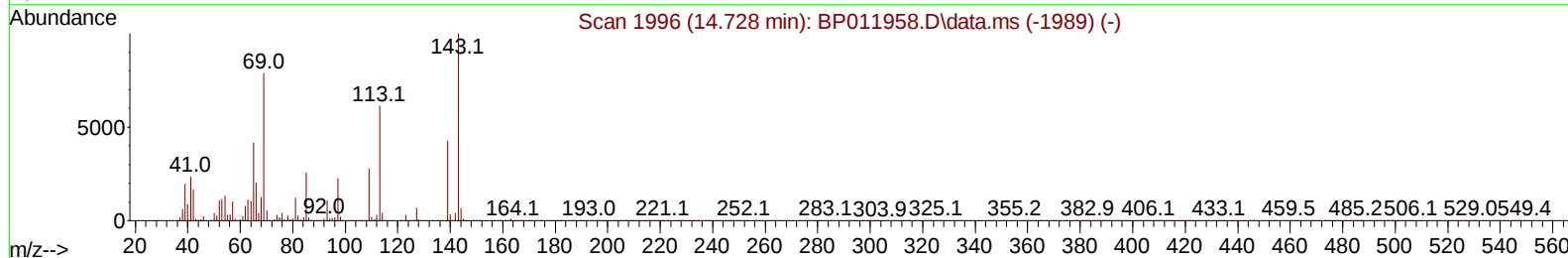
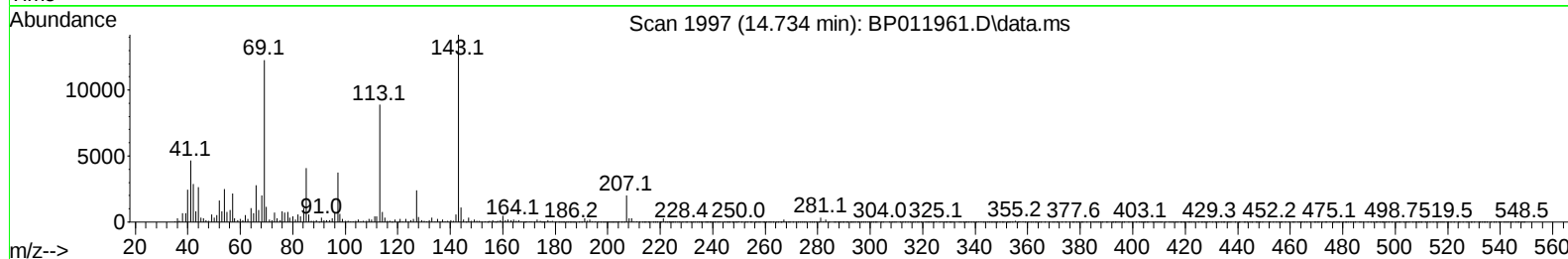
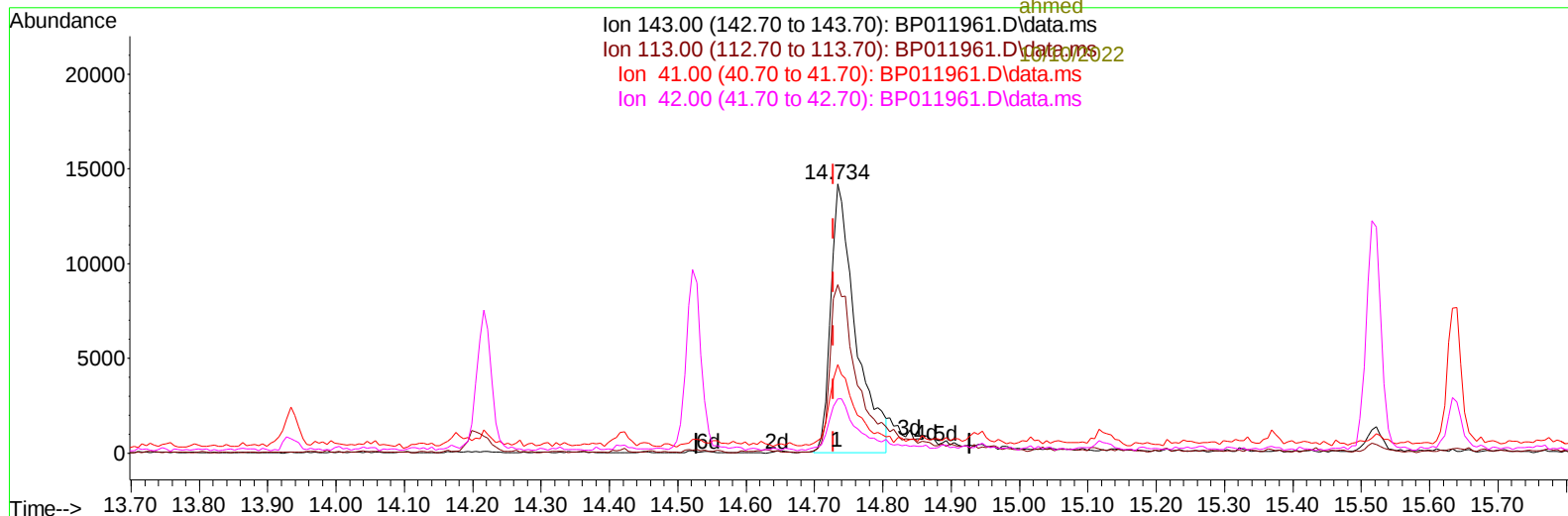
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TIC: BP011961.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 3.16 ng/ul m

response 35809

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	62.70
41.00	24.20	32.80#
42.00	16.40	20.21#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----10/10/2022						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	291866	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1238272	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	766254	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1620462	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1558246	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1558006	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	28329	4.281	ng/uL	0.00
4) Pyridine-d5	3.728	84	80440	4.150	ng/ul	0.00
7) Phenol-d5	7.040	99	112207	4.453	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	287193	20.563	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	317689	15.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	202414	9.718	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	195777	20.757	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	187951	18.626	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	354881	19.013	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	366252	12.871	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1277943	23.322	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1351749	21.640	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	35809m	3.156	ng/ul	0.00
60) Fluorene-d10	15.522	176	1154014	24.251	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	163601	16.680	ng/ul	0.00
73) Anthracene-d10	17.381	188	2008957	27.216	ng/ul	0.00
81) Pyrene-d10	19.616	212	2357071	29.868	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2261363	28.995	ng/ul	0.00

Target Compounds Qvalue

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

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