

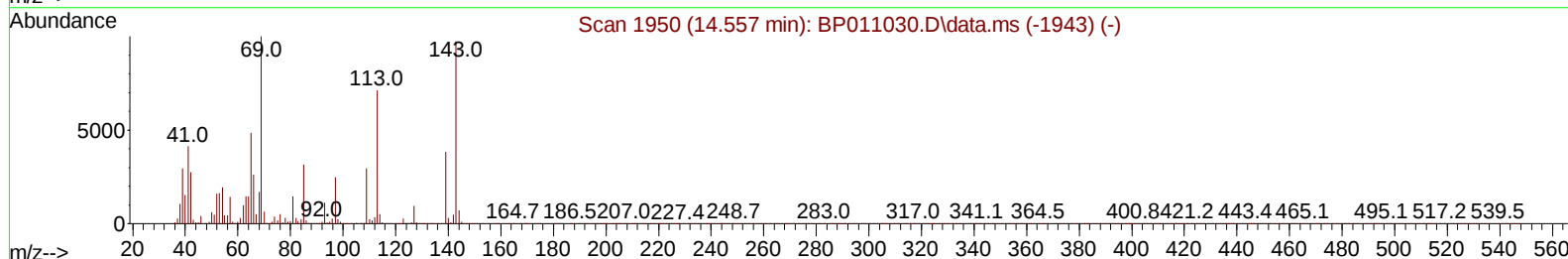
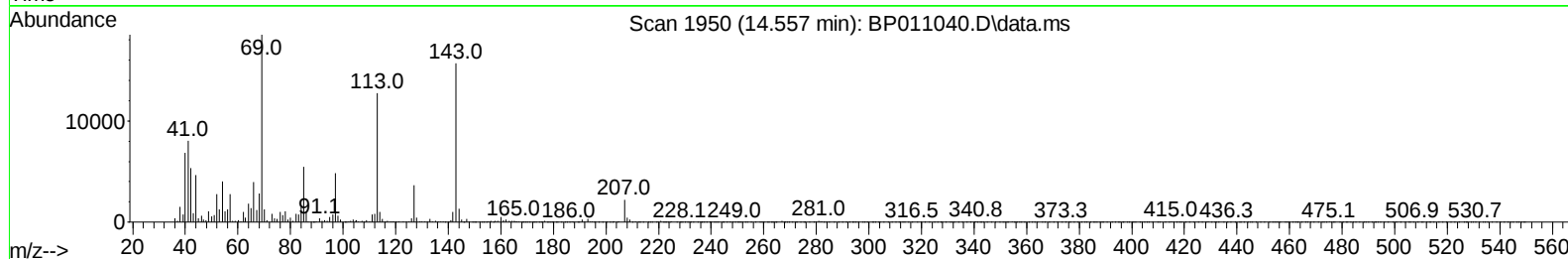
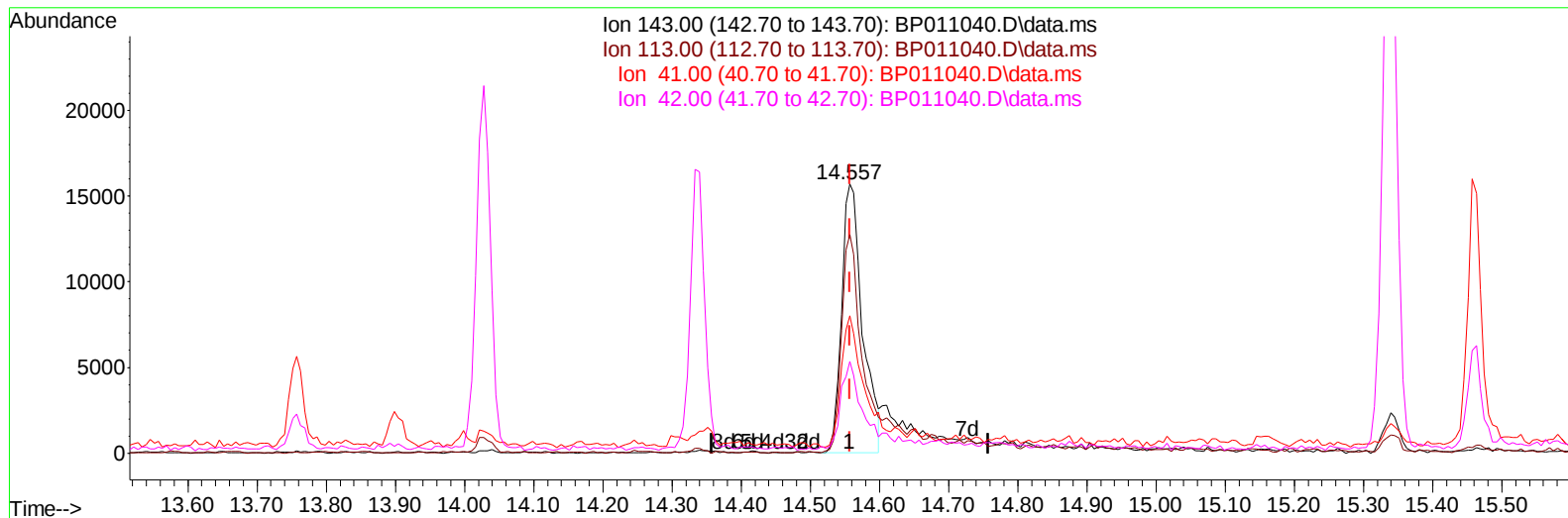
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011040.D
Acq On : 19 Jul 2022 19:34
Operator : CG/JU
Sample : N3753-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHQ4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/20/2022
Supervised By :mohammad ahmed 07/21/2022

Quant Time: Jul 19 23:41:37 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 19 03:25:54 2022
Response via : Initial Calibration



TIC: BP011040.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.557min (-0.000) 3.31 ng/u1****response 32958**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	81.35
41.00	59.00	51.02
42.00	41.10	34.04

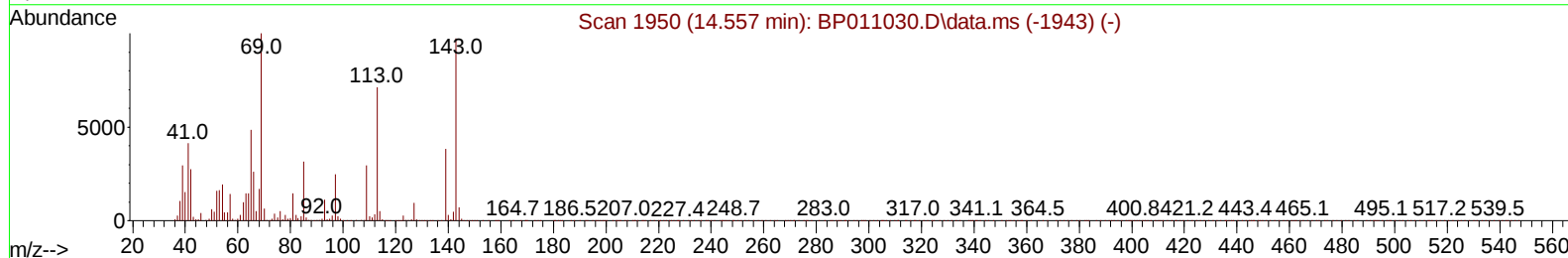
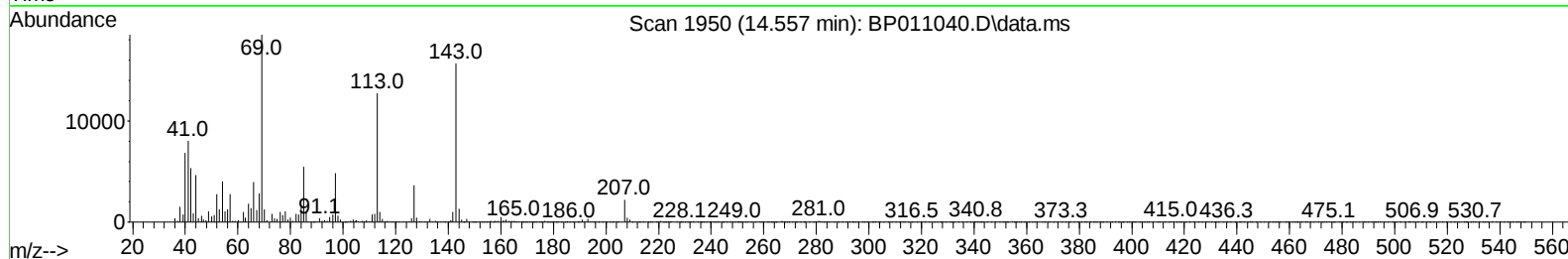
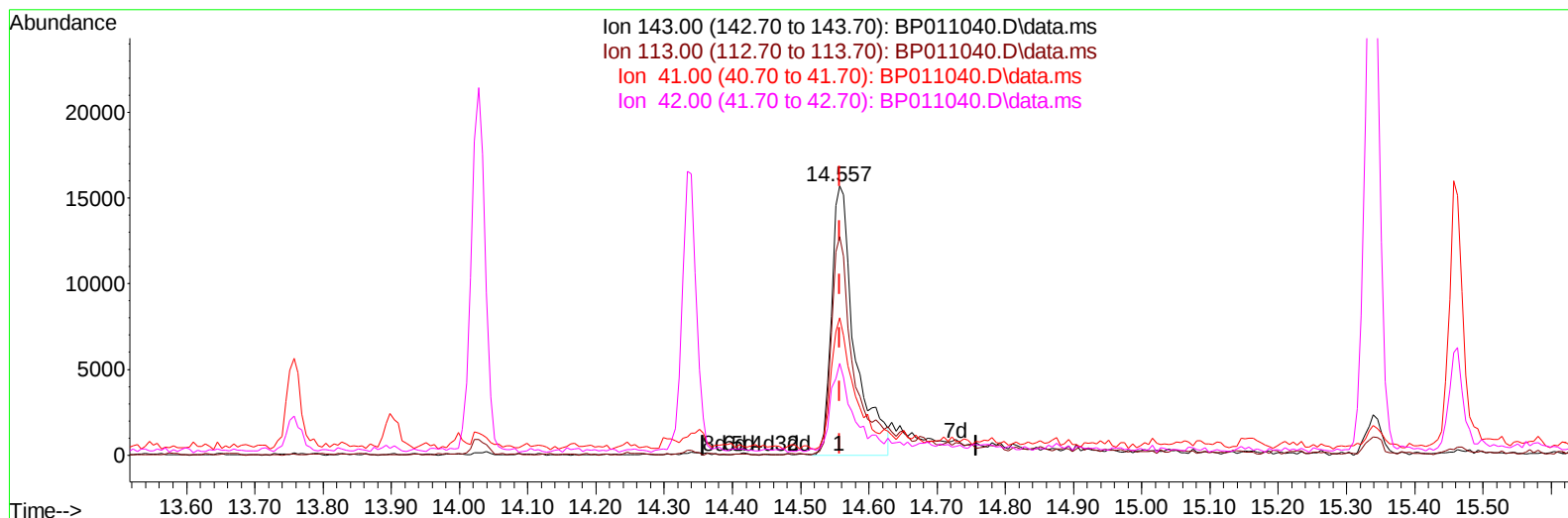
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TIC: BP011040.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.000) 3.71 ng/ul m

response 36978

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	81.35
41.00	59.00	51.02
42.00	41.10	34.04

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	330499	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	1319140	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.339	164	719674	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1439889	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1433317	20.000	ng/ul	# 0.00
88) Perylene-d12	23.580	264	1466027	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	44231	4.825	ng/uL	0.00
4) Pyridine-d5	3.593	84	229927	9.703	ng/ul	0.00
7) Phenol-d5	6.858	99	188528	6.949	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	615202	35.265	ng/ul	0.00
11) 2-Chlorophenol-d4	7.210	132	582670	26.580	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	350268	16.127	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	352695	36.688	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	310803	32.674	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	529933	30.463	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	786227	27.415	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	1997114	40.602	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160	2335374	40.241	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	36978m	3.715	ng/ul	0.00
60) Fluorene-d10	15.339	176	1724167	41.239	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	191623	26.894	ng/ul	0.00
73) Anthracene-d10	17.204	188	2872777	46.931	ng/ul	0.00
81) Pyrene-d10	19.463	212	3306912	43.929	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.433	264	3289367	46.716	ng/ul	0.00

Target Compounds Qvalue

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

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