

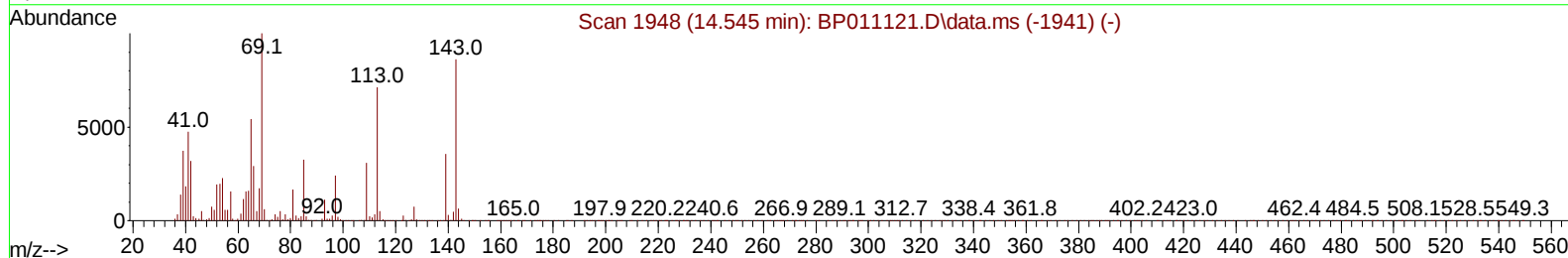
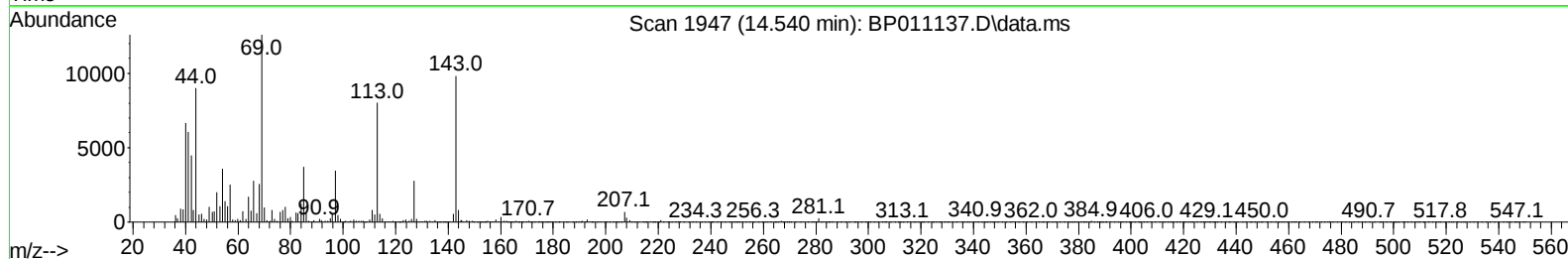
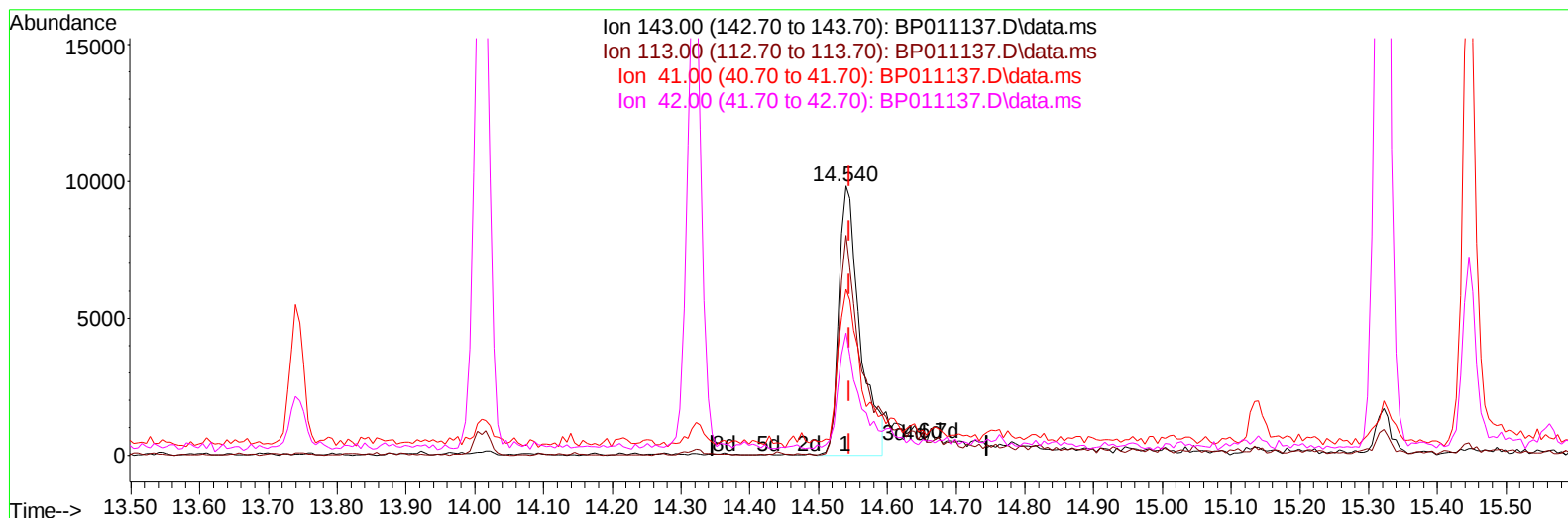
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011137.D
 Acq On : 27 Jul 2022 19:24
 Operator : CG/JU
 Sample : N3871-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJ13

Manual IntegrationsAPPROVED

Quant Time: Jul 28 00:03:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/28/2022
 Supervised By :mohammad ahmed 07/28/2022



TIC: BP011137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (-0.006) 2.65 ng/u1

response 21184

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	81.63
41.00	59.00	61.63
42.00	41.10	45.46

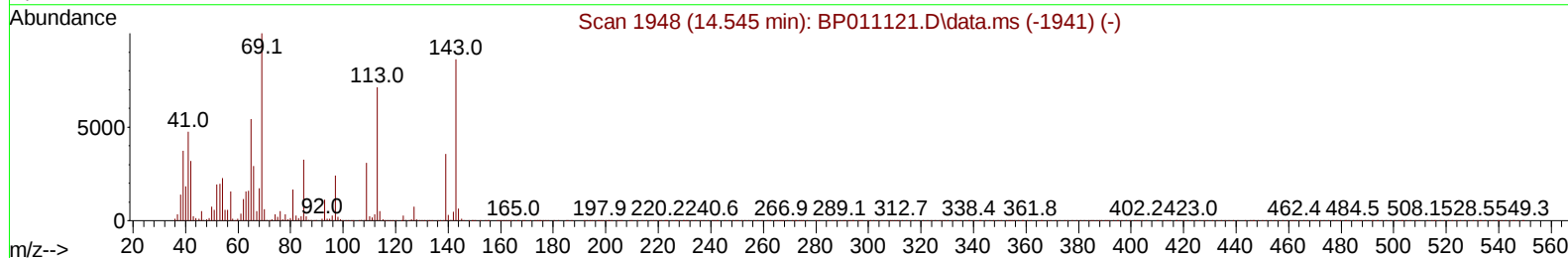
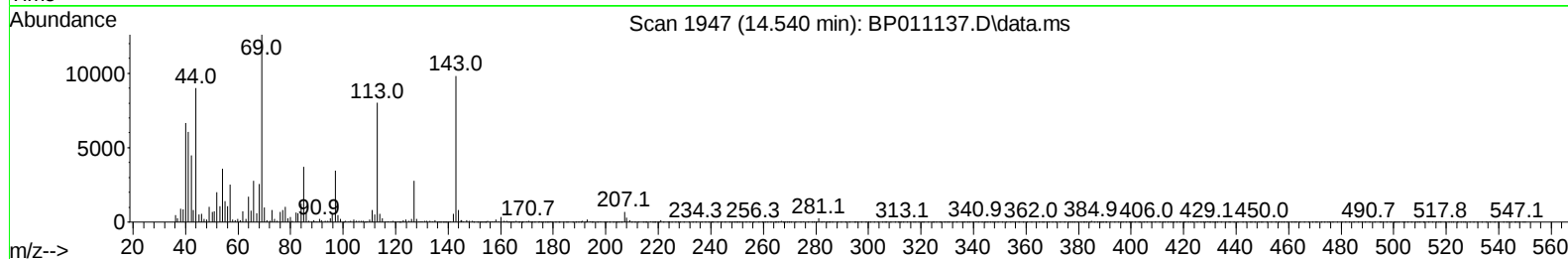
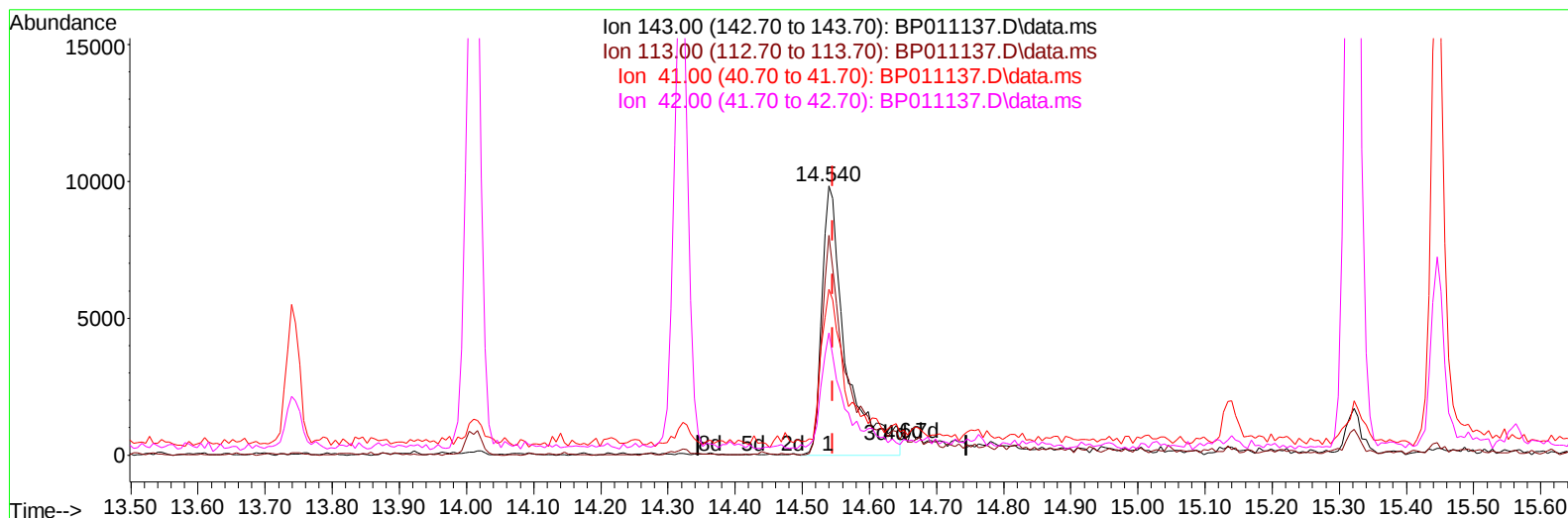
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TIC: BP011137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (-0.006) 3.07 ng/ul m

response 24548

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	81.63
41.00	59.00	61.63
42.00	41.10	45.46

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.658	152	280563	20.000	ng/ul	0.00
20) Naphthalene-d8	10.452	136	1139587	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.322	164	577157	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	1101721	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.204	240	1114842	20.000	ng/ul	#-0.01
88) Perylene-d12	23.545	264	1207454	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	32464	4.171	ng/uL	0.00
4) Pyridine-d5	3.575	84	171929	8.547	ng/ul	0.00
7) Phenol-d5	6.840	99	147763	6.416	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	564372	38.109	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	476936	25.629	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	279796	15.176	ng/ul	0.00
21) Nitrobenzene-d5	8.822	128	311512	37.510	ng/ul	0.00
24) 2-Nitrophenol-d4	9.540	143	281253	34.226	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	440047	29.282	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	687420	27.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.745	166	1497557	37.964	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160	1872995	40.243	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	24548m	3.075	ng/ul	0.00
60) Fluorene-d10	15.322	176	1288513	38.429	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	163687	30.025	ng/ul	-0.01
73) Anthracene-d10	17.186	188	2048296	43.733	ng/ul	0.00
81) Pyrene-d10	19.445	212	2433517	41.562	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.392	264	2550237	43.975	ng/ul	-0.01

Target Compounds Qvalue

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

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