

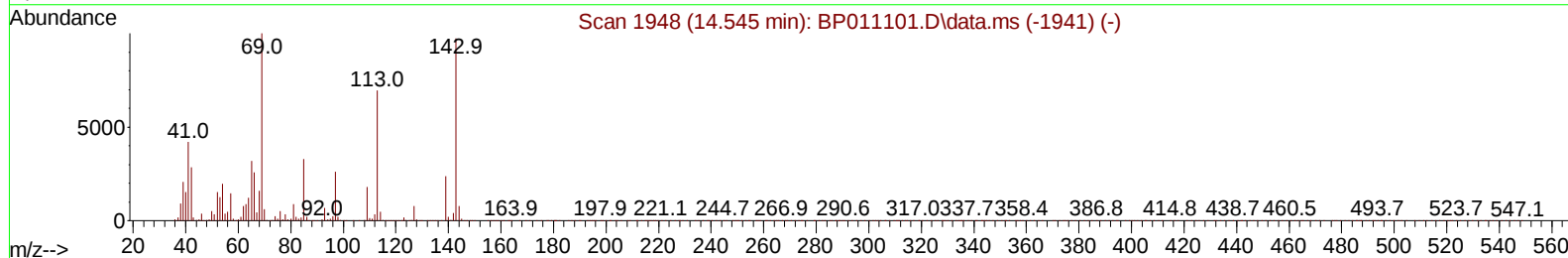
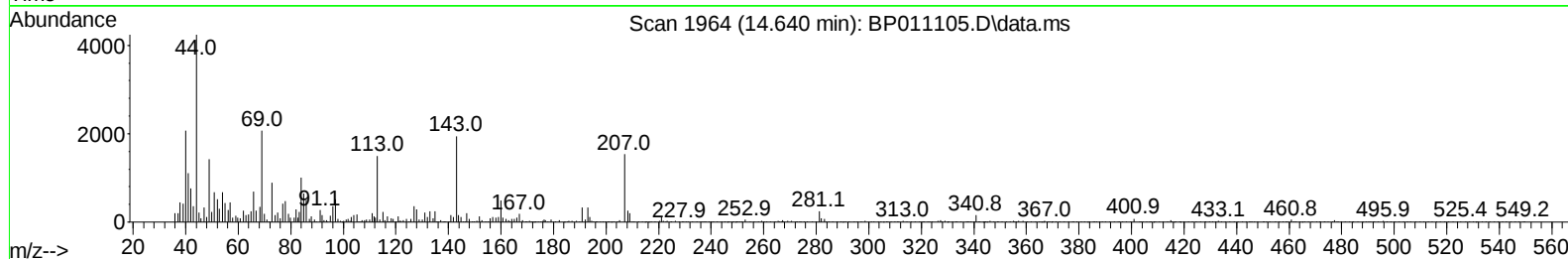
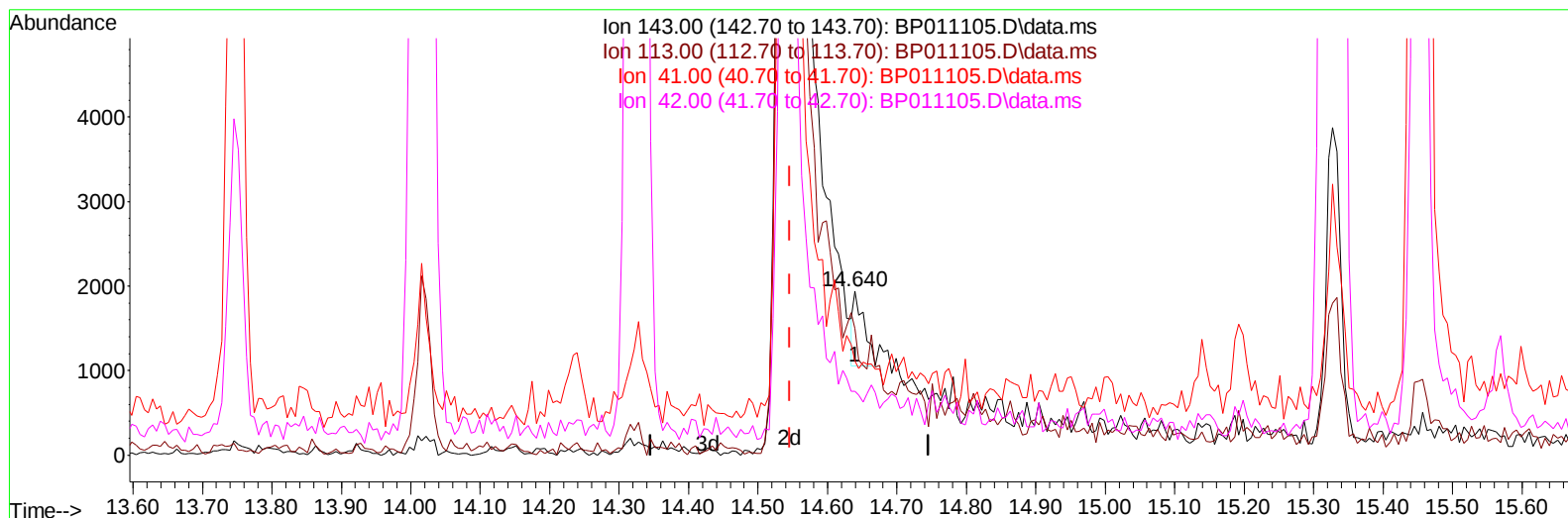
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011105.D
 Acq On : 26 Jul 2022 12:11
 Operator : CG/JU
 Sample : N3801-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHX0

Manual IntegrationsAPPROVED

Quant Time: Jul 26 14:43:09 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 07/27/2022
 Supervised By :mohammad ahmed 07/27/2022



TIC: BP011105.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.640min (+ 0.094) 0.05 ng/u1

response 943

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	76.94
41.00	59.00	57.18
42.00	41.10	38.96

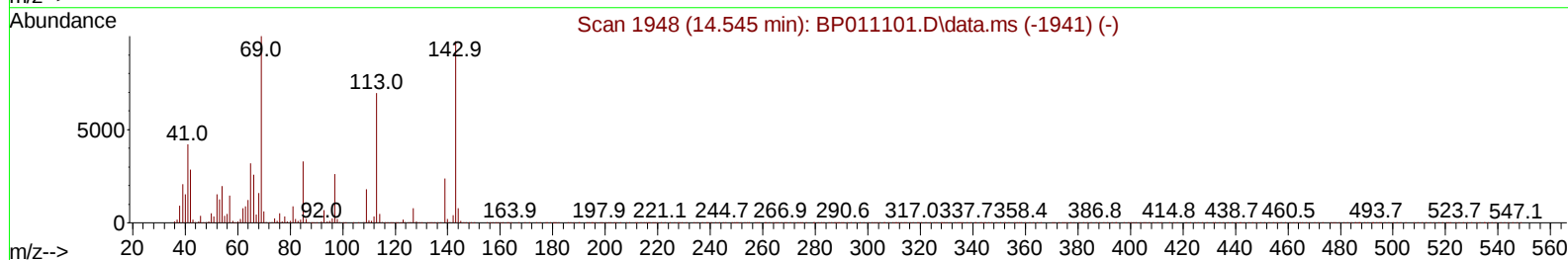
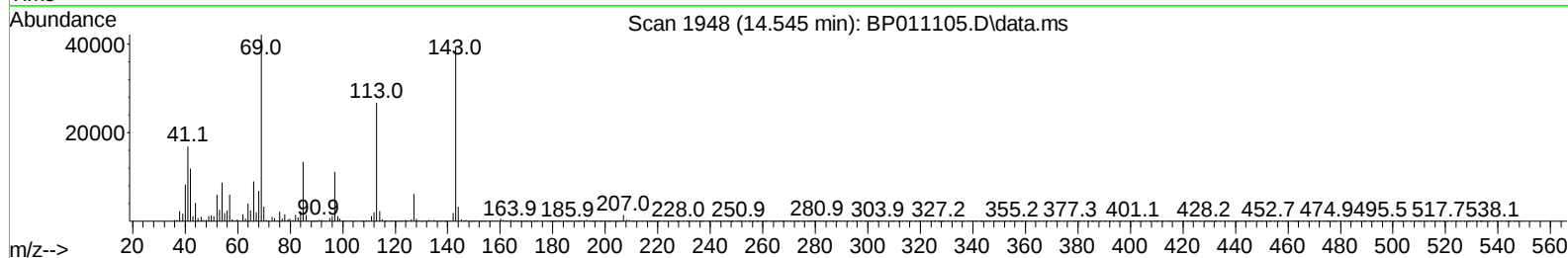
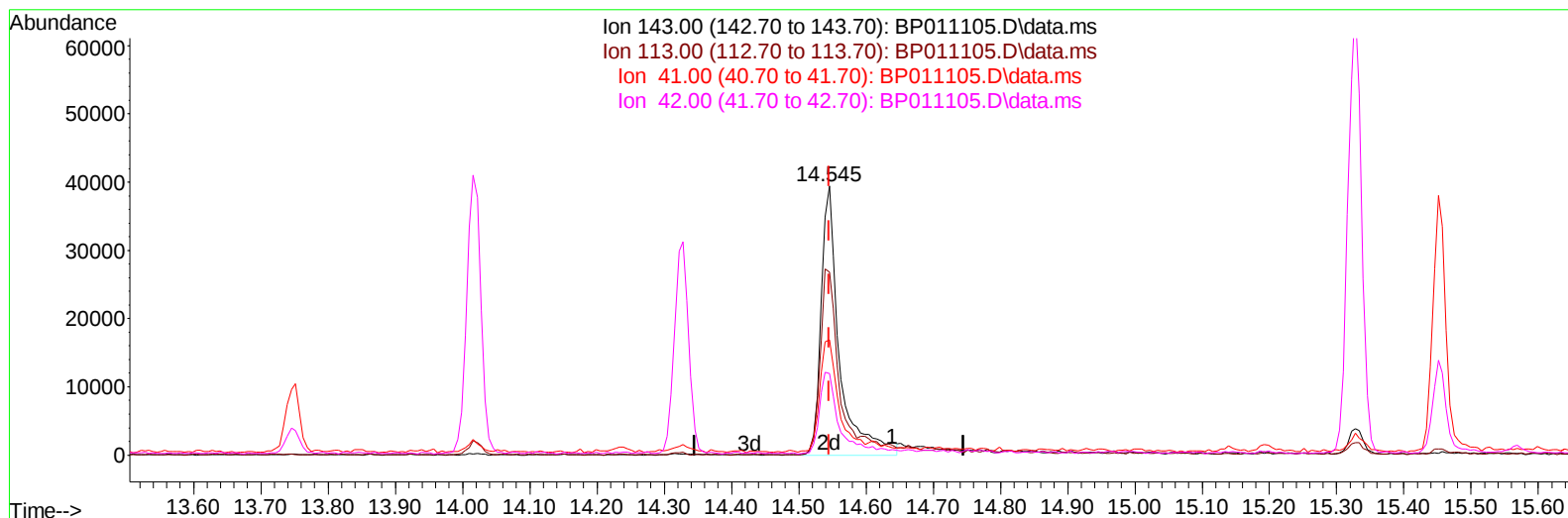
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(54) 4-Nitrophenol-d4 (S)

14.545min (0.000) 4.12 ng/ul m

response 74718

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	68.01#
41.00	59.00	42.87#
42.00	41.10	30.30#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	499203	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	2247598	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.328	164	1311514	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	2300832	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	1426149	20.000	ng/ul	# 0.00
88) Perylene-d12	23.551	264	1464808	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	67767	4.894	ng/uL	0.00
4) Pyridine-d5	3.575	84	289931	8.101	ng/ul	0.00
7) Phenol-d5	6.840	99	311518	7.602	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	1023188	38.831	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	935946	28.267	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	618648	18.858	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	643214	39.270	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	645024	39.798	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	987877	33.329	ng/ul	0.00
31) 4-Chloroaniline-d4	10.604	131	1472642	30.138	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	3907477	43.592	ng/ul	0.00
49) Acenaphthylene-d8	14.016	160	4539441	42.922	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	74718m	4.119	ng/ul	0.00
60) Fluorene-d10	15.328	176	3119207	40.939	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	457641	40.196	ng/ul	0.00
73) Anthracene-d10	17.192	188	4307718	44.040	ng/ul	0.00
81) Pyrene-d10	19.445	212	4010095	53.538	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	3165511	44.994	ng/ul	0.00

Target Compounds Qvalue

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

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