

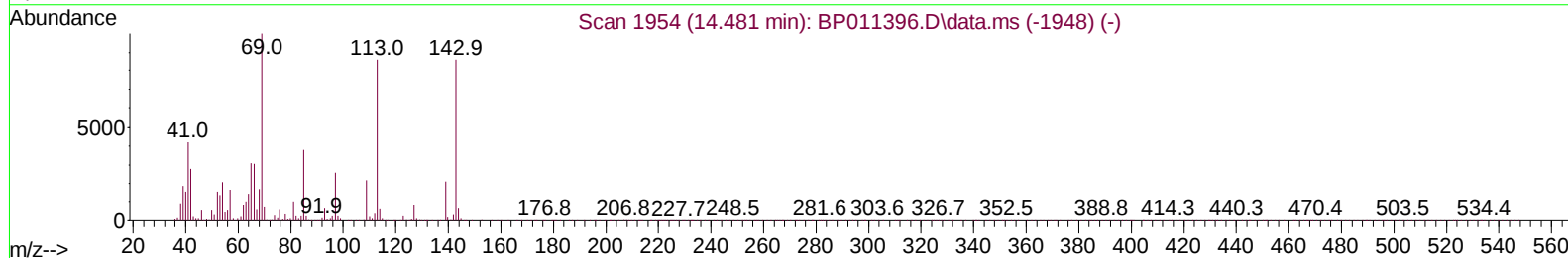
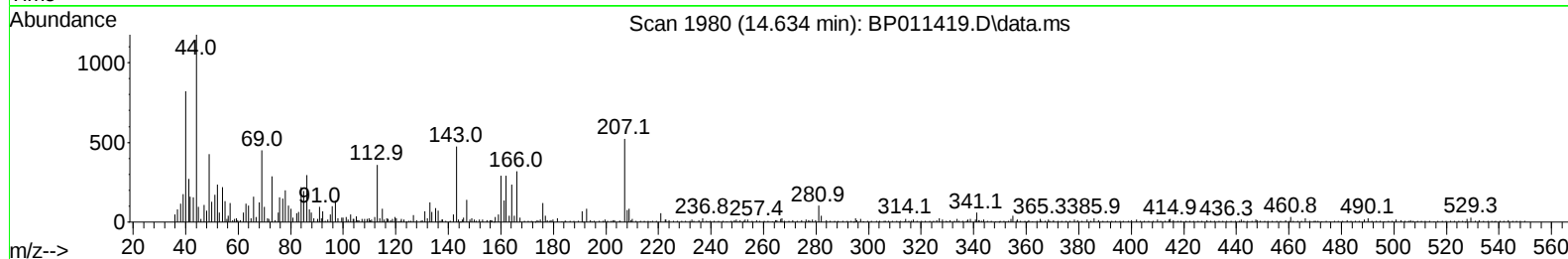
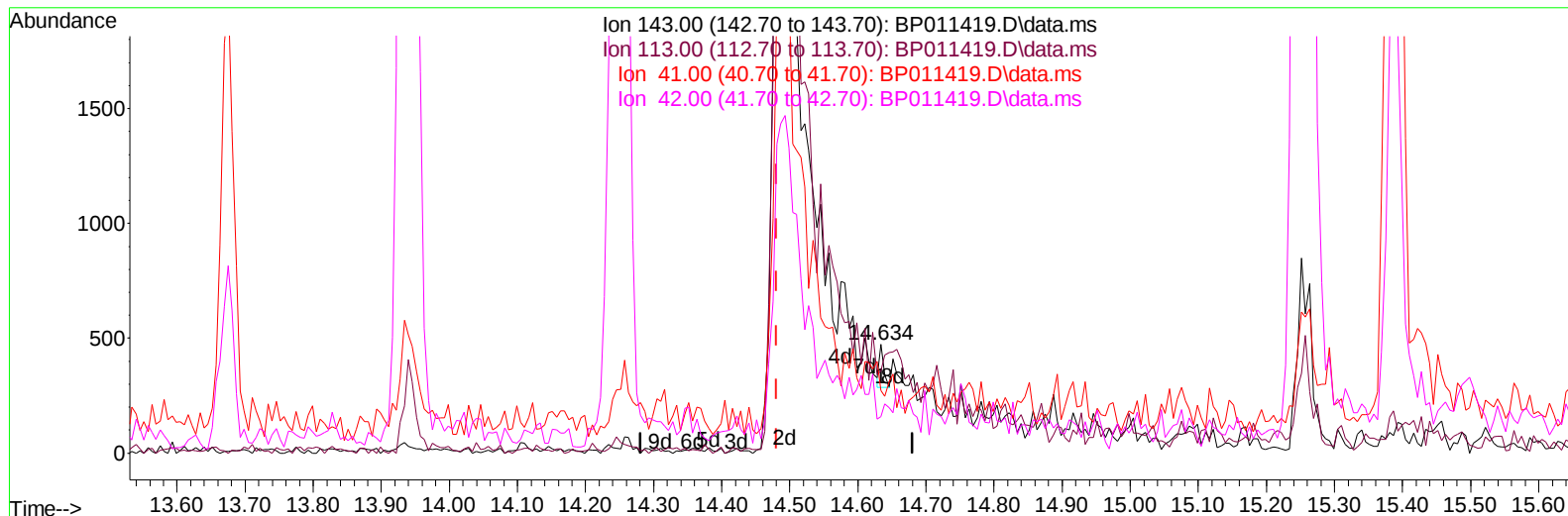
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011419.D
Acq On : 12 Aug 2022 11:47
Operator : CG/JU
Sample : N4129-07
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGBT0

Manual IntegrationsAPPROVED

Quant Time: Aug 12 12:24:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011419.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.153) 0.04 ng/u1

response 86

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	75.90
41.00	54.40	57.51
42.00	35.50	33.62

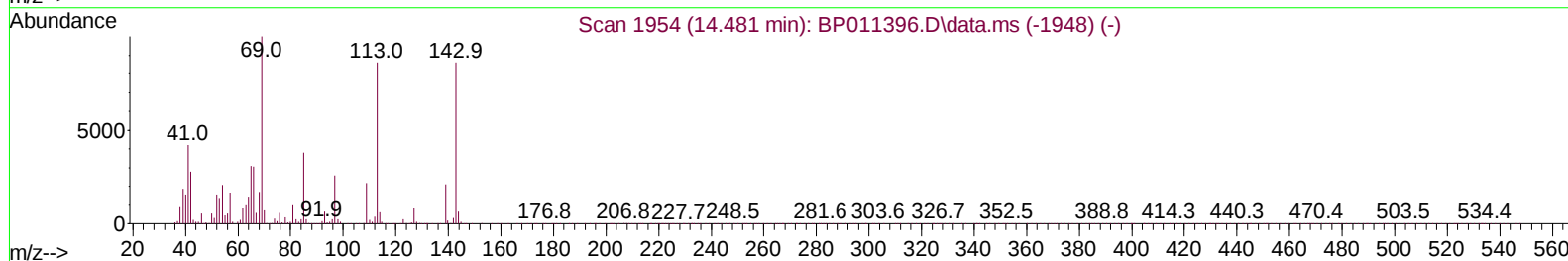
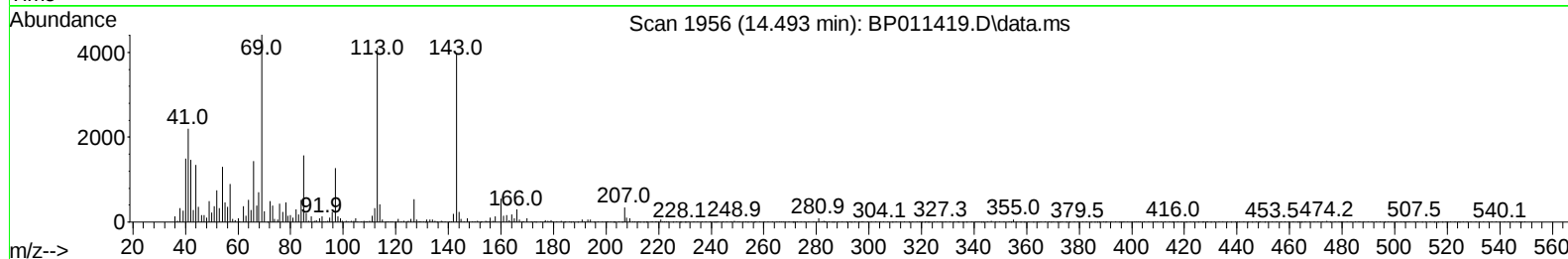
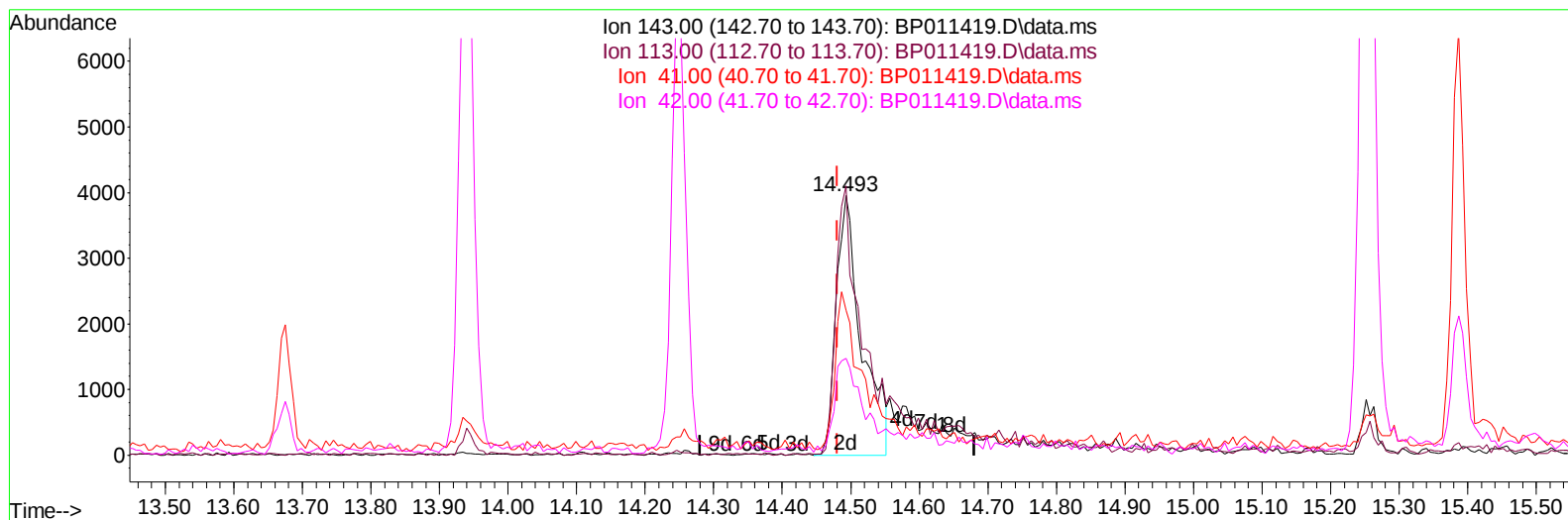
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TIC: BP011419.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.012) 4.23 ng/ul m

response 10182

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	103.64#
41.00	54.40	55.82
42.00	35.50	37.20

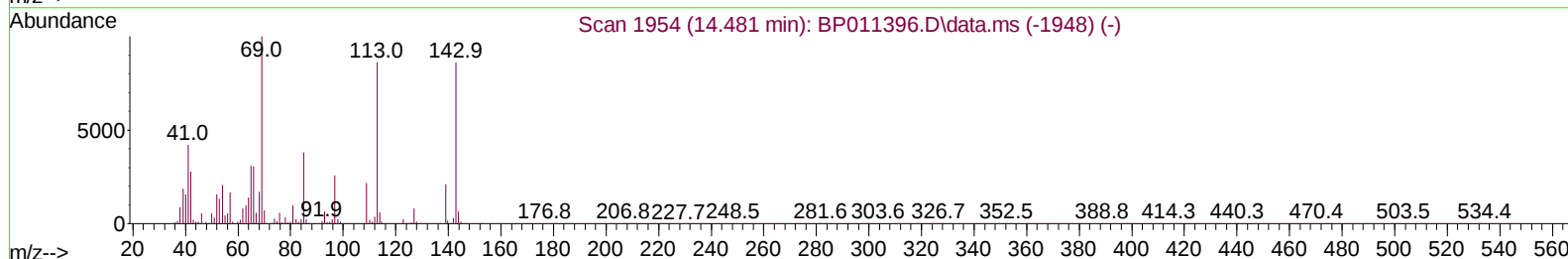
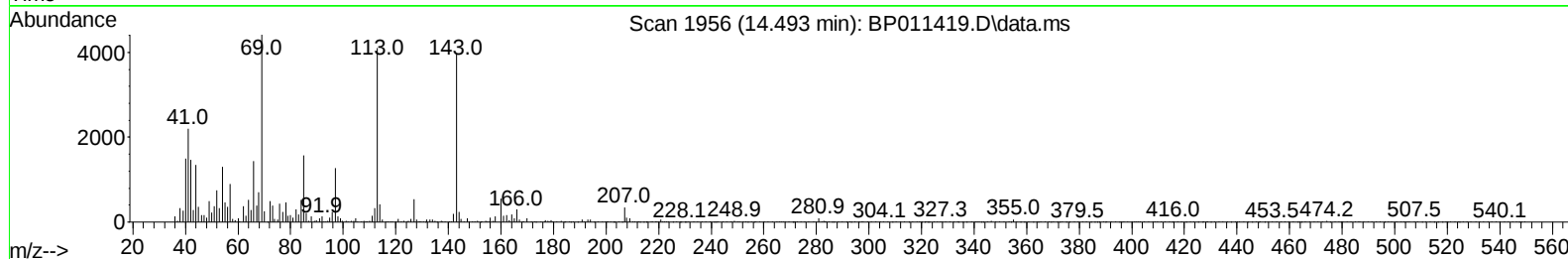
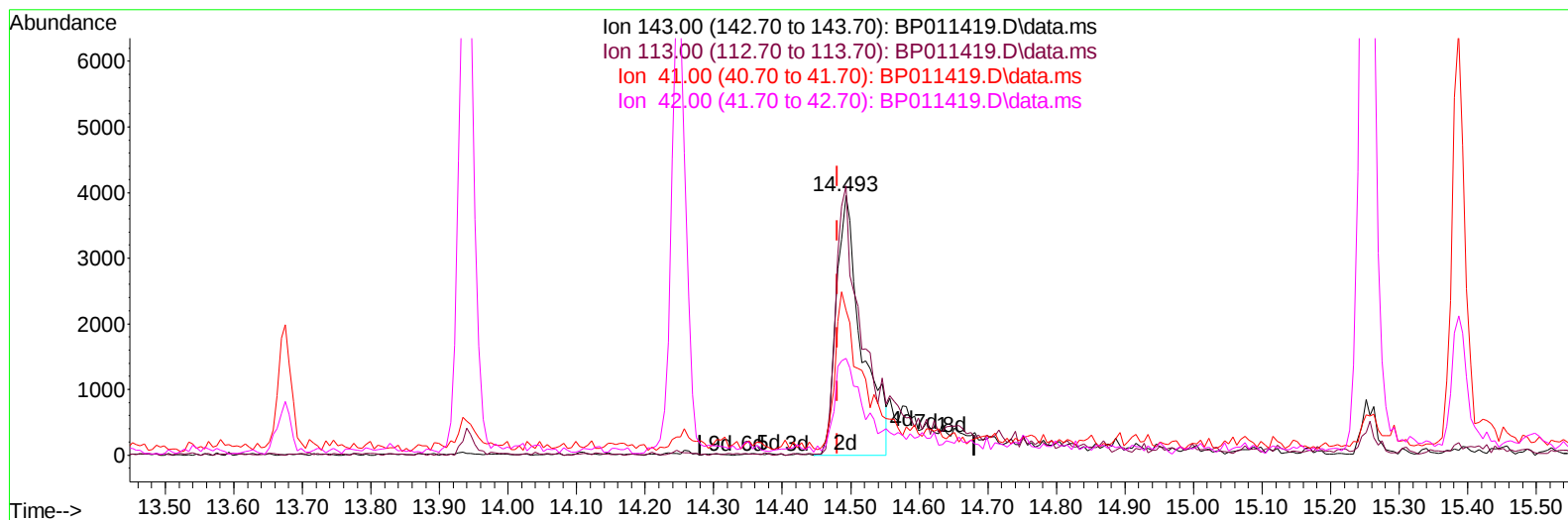
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	77631	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	328545	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	189805	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	391599	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	381359	20.000	ng/ul	# 0.00
88) Perylene-d12	23.463	264	398589	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	12676	5.006	ng/uL	0.00
4) Pyridine-d5	3.499	84	49273	7.714	ng/ul	0.01
7) Phenol-d5	6.758	99	49473	7.201	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	148647	30.515	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	138468	25.688	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	88899	16.480	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	86965	38.893	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	79817	38.989	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	125588	30.378	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	179602	24.850	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	532978	40.470	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	582613	36.751	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	10182m	4.233	ng/ul	0.01
60) Fluorene-d10	15.257	176	431035	38.271	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	65468	35.993	ng/ul	0.00
73) Anthracene-d10	17.122	188	767650	43.613	ng/ul	0.00
81) Pyrene-d10	19.386	212	891029	42.409	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	873314	43.446	ng/ul	0.00

Target Compounds Qvalue

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

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