

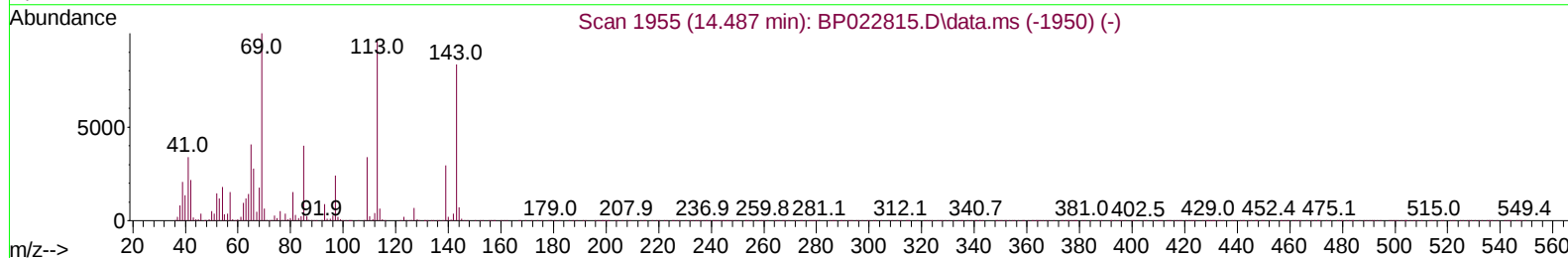
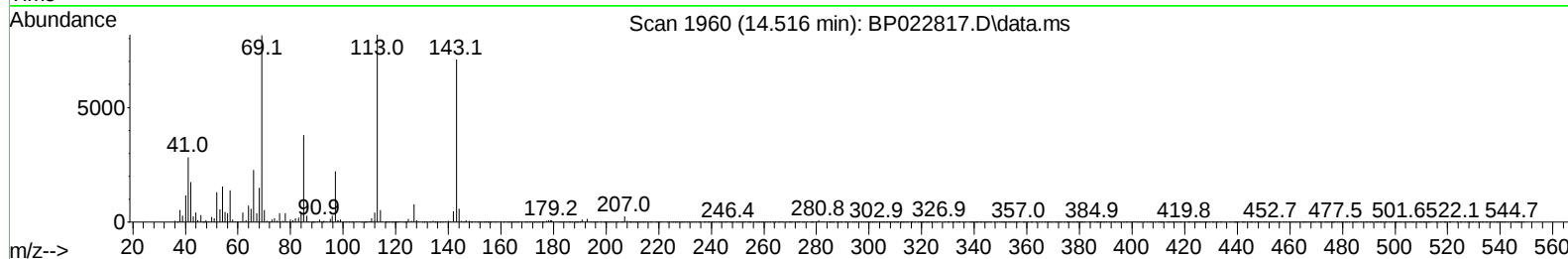
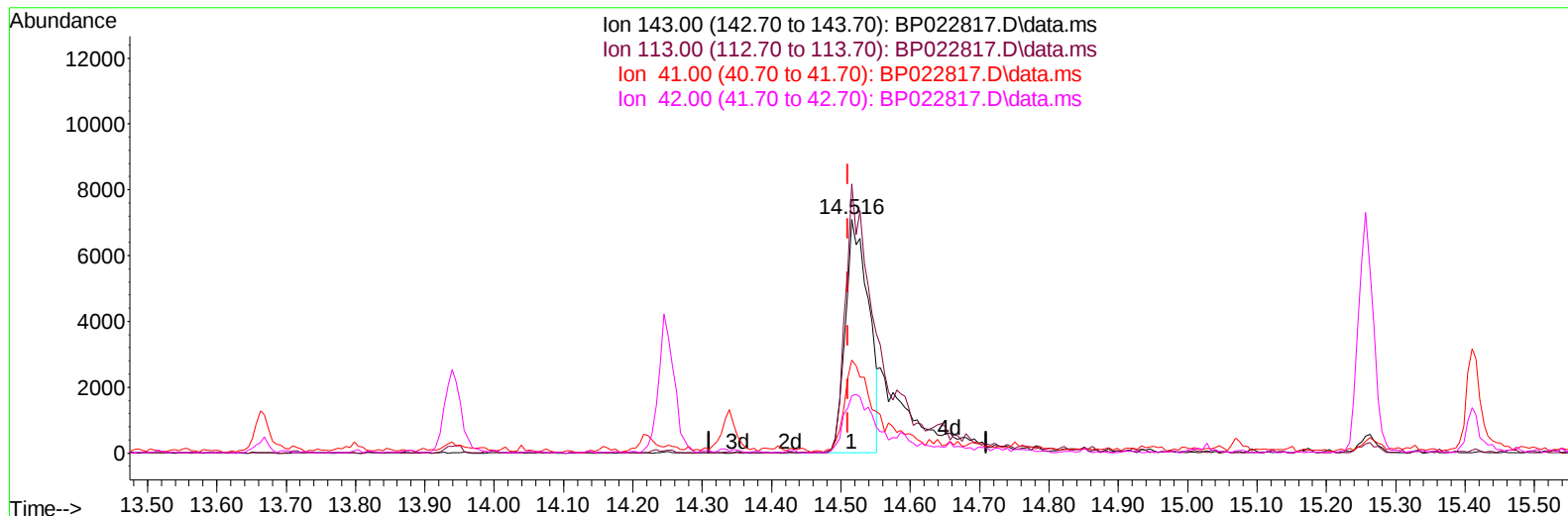
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022817.D
Acq On : 02 Nov 2024 00:17
Operator : RC/JU
Sample : P4630-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP13

Manual IntegrationsAPPROVED

Quant Time: Nov 04 05:10:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022817.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 6.84 ng/u1

response 16396

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	115.08
41.00	35.90	39.90
42.00	23.70	24.51

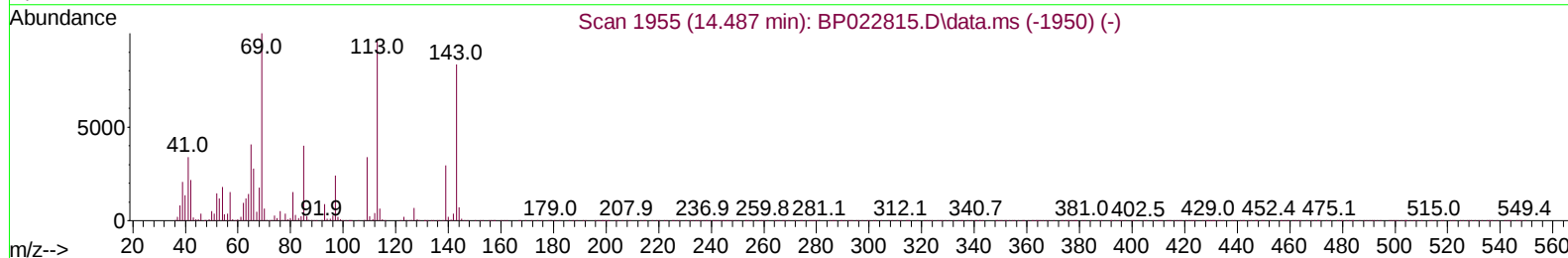
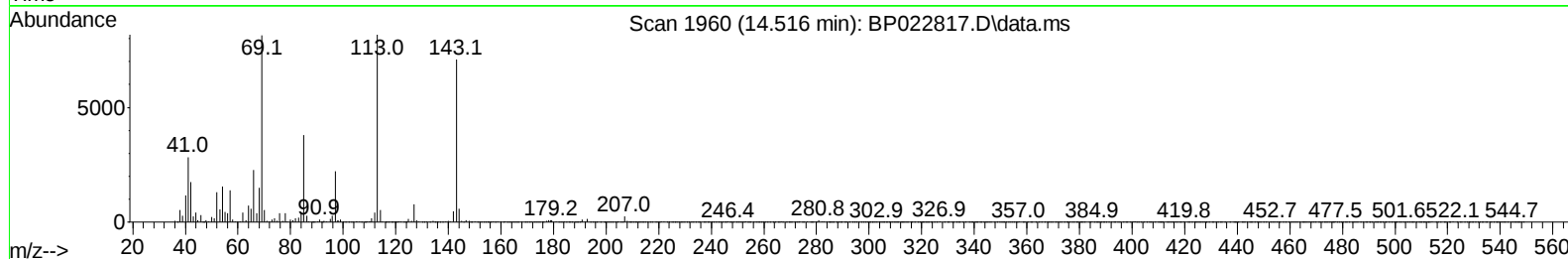
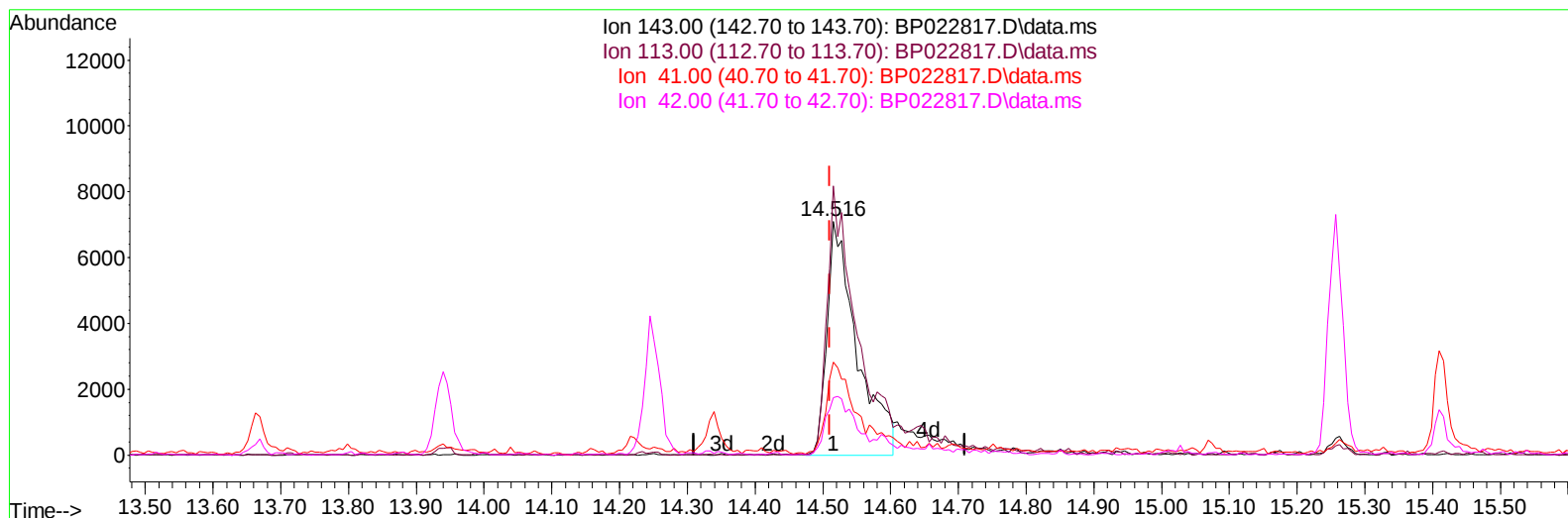
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(54) 4-Nitrophenol-d4 (S)

14.516min (+ 0.006) 9.08 ng/ul m

response 21778

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	115.08
41.00	35.90	39.90
42.00	23.70	24.51

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Quant Time: Nov 04 05:11:39 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	61781	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.386	136	224498	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	179928	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.051	188	451673	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.498	240	520694	20.000	ng/ul	# 0.00
88) Perylene-d12	24.762	264	640772	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5642	3.549	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.828	99	41535	7.987	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	84541	27.659	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	106971	28.988	ng/ul	-0.01
15) 4-Methylphenol-d8	8.340	113	78841	18.996	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	54789	31.740	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	65769	32.910	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	131005	30.843	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.669	166	473046	33.090	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	310155	20.427	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.516	143	21778m	9.081	ng/ul	0.00
60) Fluorene-d10	15.257	176	407021	32.825	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.410	200	71849	24.187	ng/ul	0.00
73) Anthracene-d10	17.163	188	608738	28.649	ng/ul	0.00
81) Pyrene-d10	19.533	212	809935	29.414	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.521	264	852765	24.920	ng/ul	0.00

Target Compounds Qvalue

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

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