

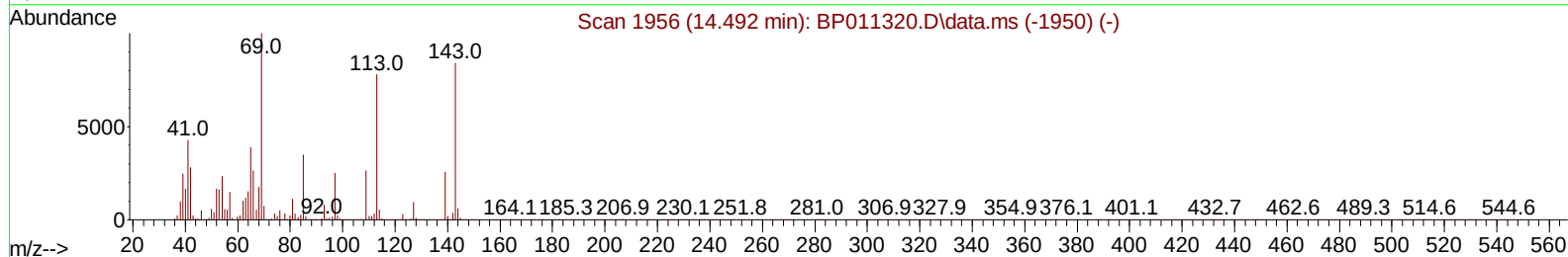
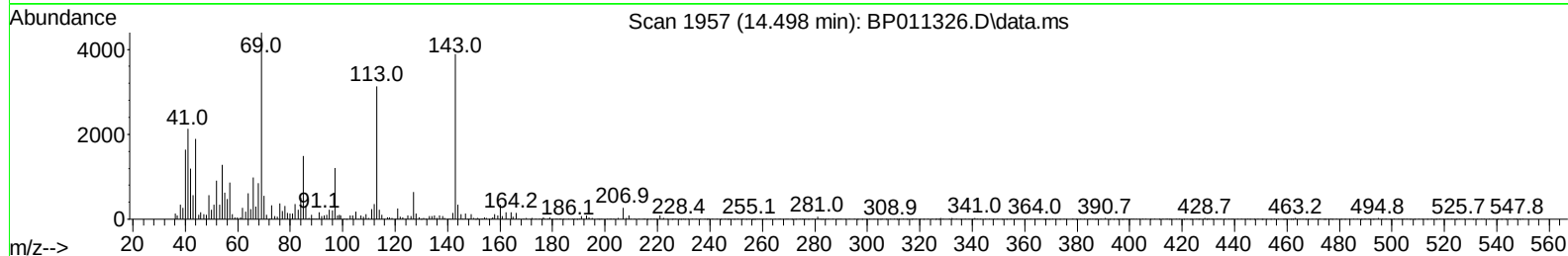
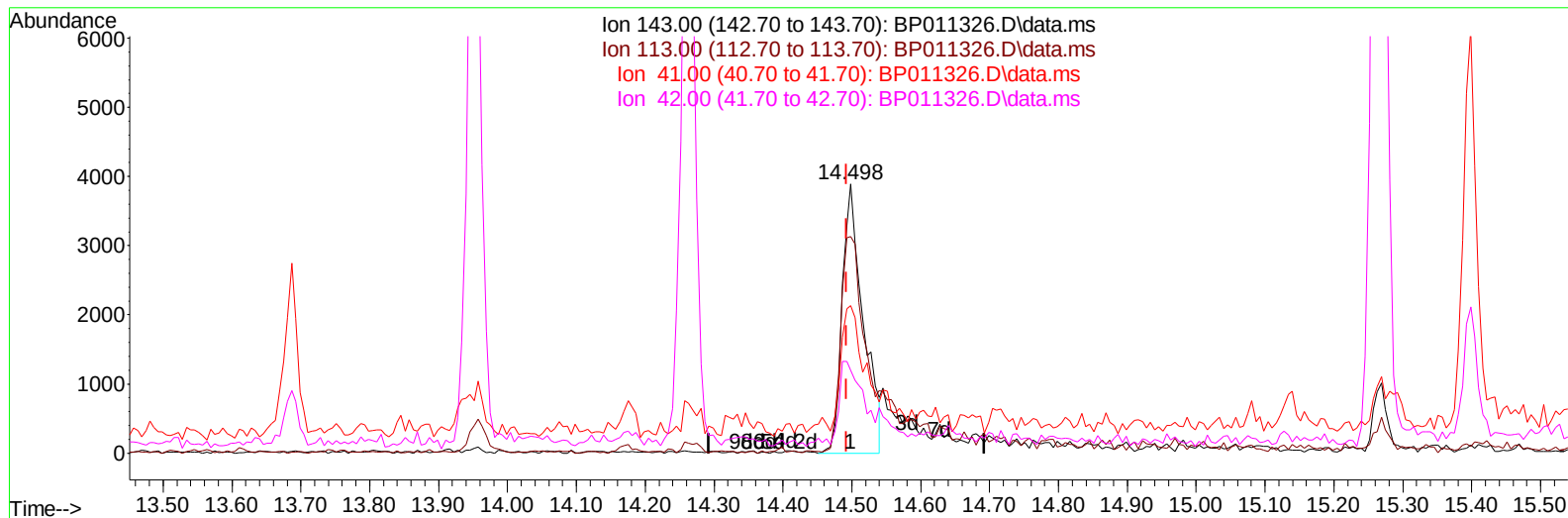
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011326.D
Acq On : 09 Aug 2022 13:10
Operator : CG/JU
Sample : N4026-22
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBJJ7

Manual IntegrationsAPPROVED

Quant Time: Aug 09 23:10:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011326.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.37 ng/ul

response 8182

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	80.59
41.00	54.40	54.77
42.00	35.50	30.80

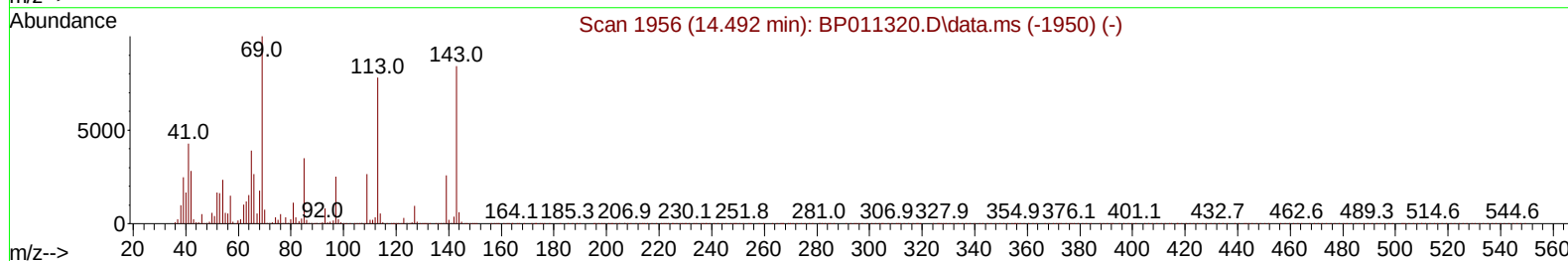
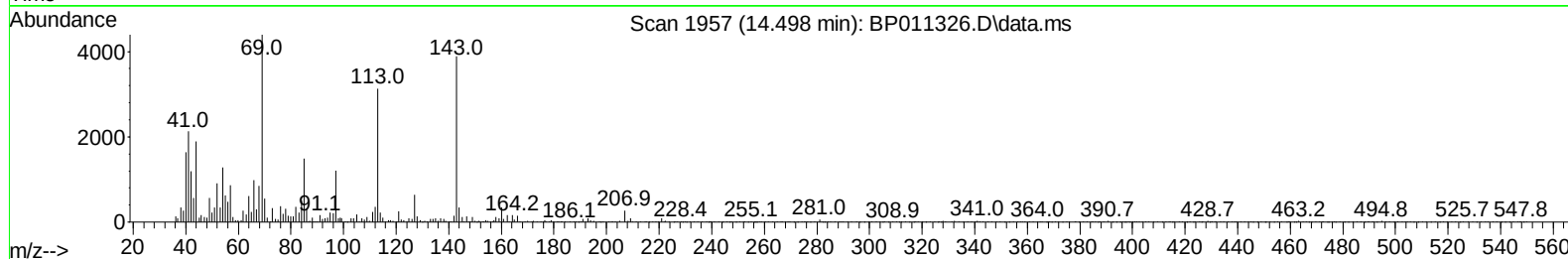
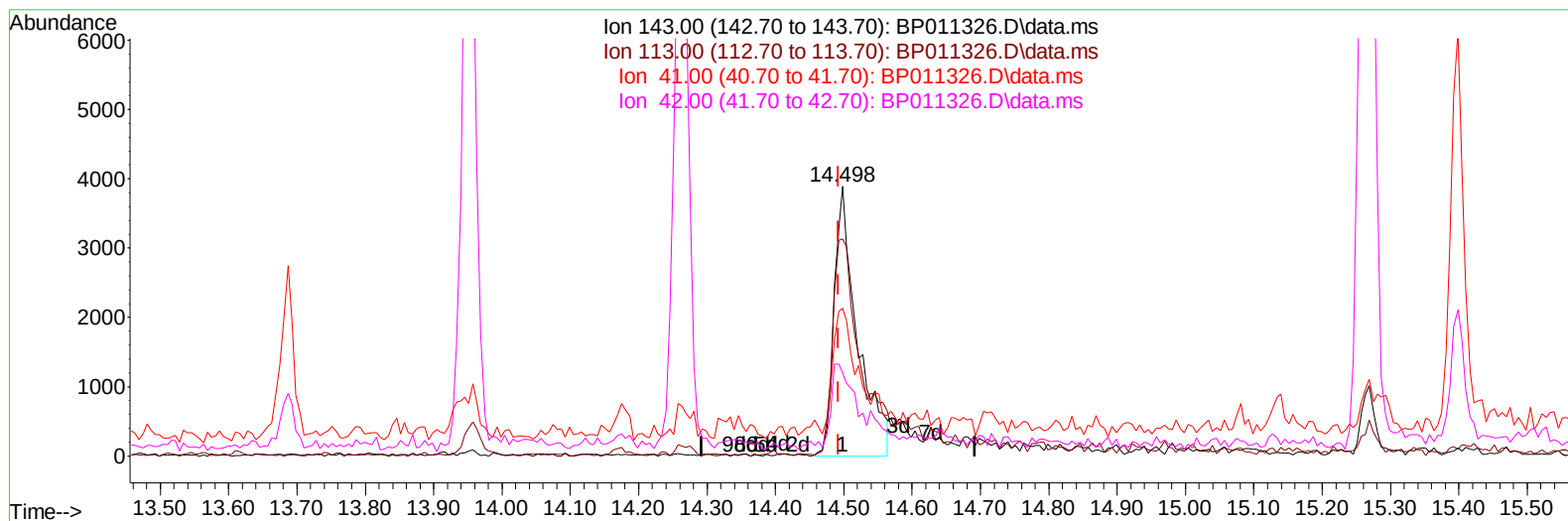
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TIC: BP011326.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.64 ng/ul m

response 9137

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	80.59
41.00	54.40	54.77
42.00	35.50	30.80

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	117604	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	495525	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	272693	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	546475	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	504217	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	509160	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	14419	3.759	ng/uL	0.01
4) Pyridine-d5	3.528	84	17735	1.833	ng/ul	0.03
7) Phenol-d5	6.775	99	56262	5.406	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	201221	27.267	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	176890	21.662	ng/ul	0.00
15) 4-Methylphenol-d8	8.311	113	104971	12.845	ng/ul	0.00
21) Nitrobenzene-d5	8.758	128	103541	30.702	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	93237	30.197	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	170140	27.287	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	109885	10.080	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	637508	33.694	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	734037	32.229	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	9137m	2.644	ng/ul	0.00
60) Fluorene-d10	15.269	176	523723	32.366	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	60339	23.771	ng/ul	0.00
73) Anthracene-d10	17.134	188	799368	32.544	ng/ul	0.00
81) Pyrene-d10	19.398	212	918718	33.072	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	835027	32.520	ng/ul	0.00

Target Compounds Qvalue

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

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