

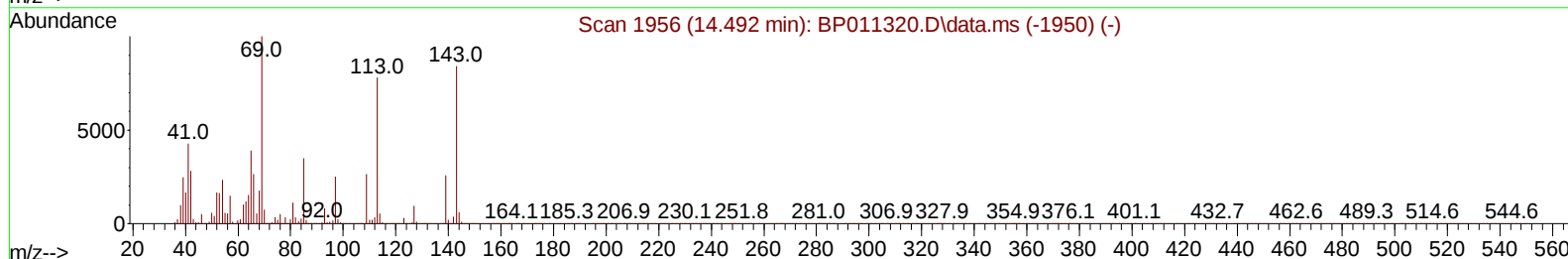
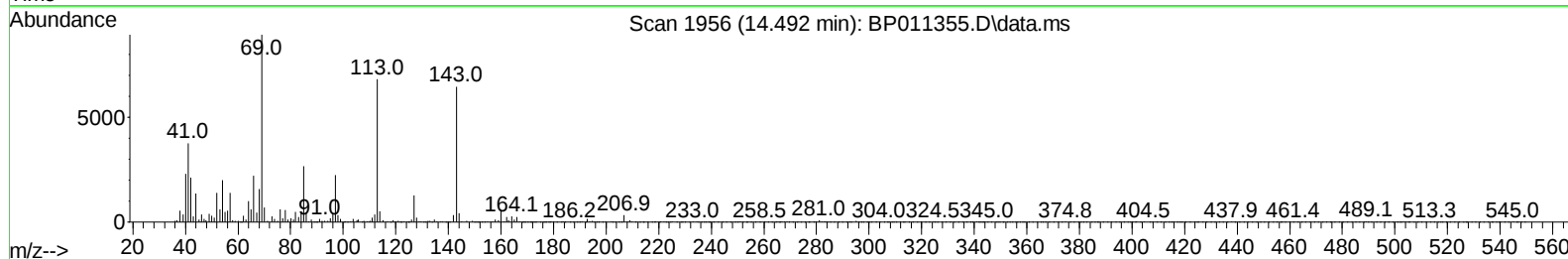
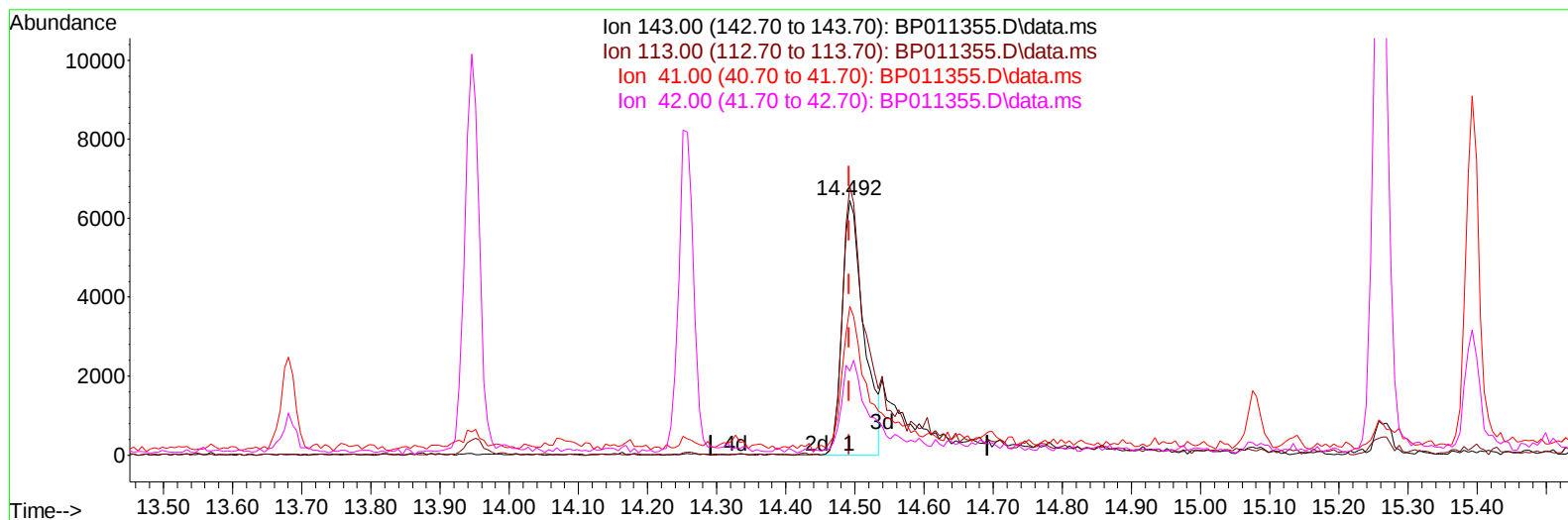
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011355.D
 Acq On : 10 Aug 2022 06:34
 Operator : CG/JU
 Sample : N4069-05
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJK3

Manual IntegrationsAPPROVED

Quant Time: Aug 10 07:00:58 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: BP011355.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.000) 4.52 ng/u1

response 13984

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	105.61#
41.00	54.40	58.36
42.00	35.50	32.92

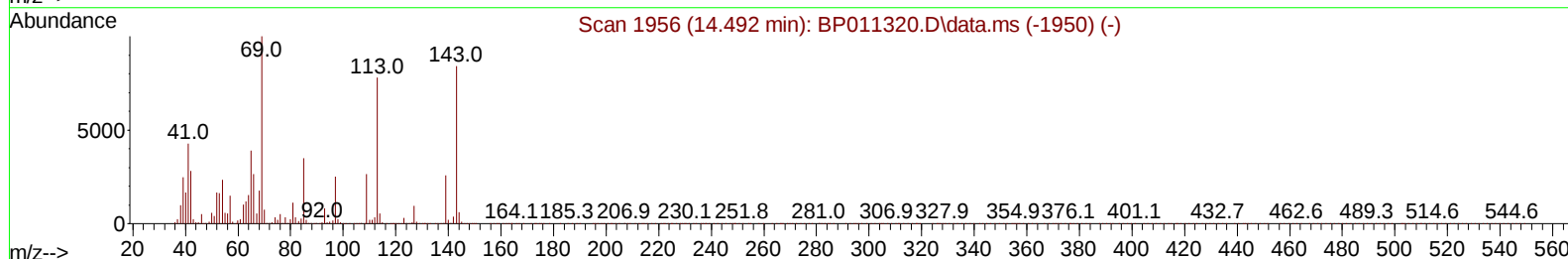
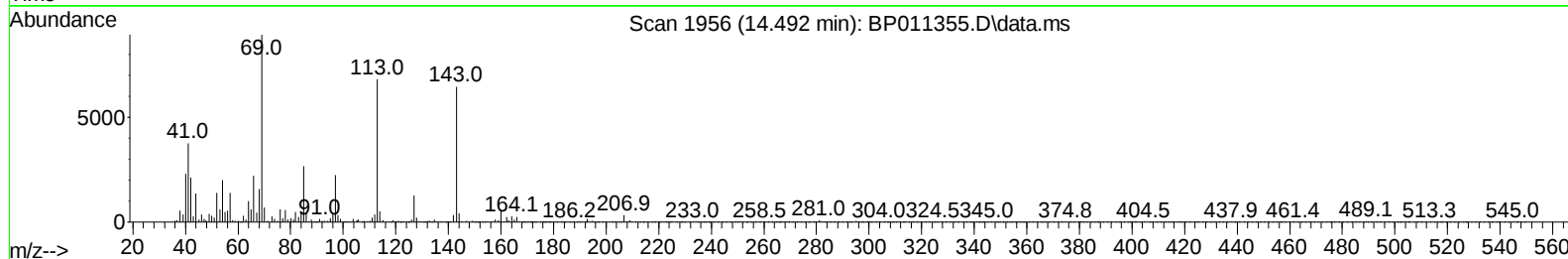
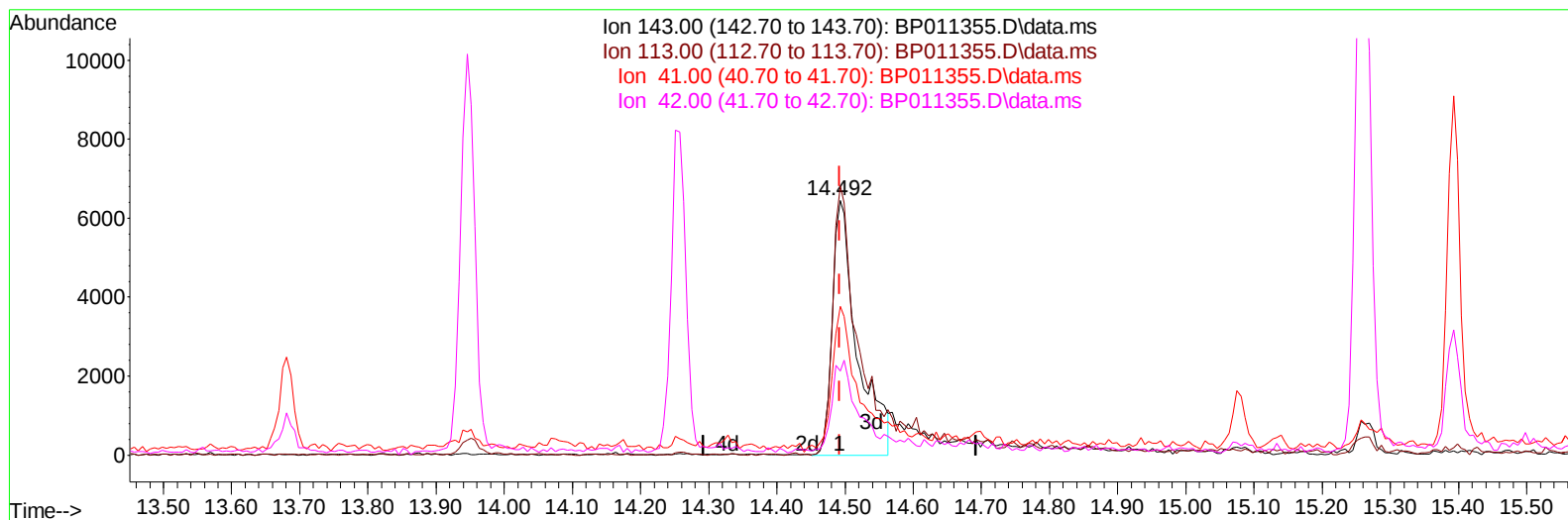
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TIC: BP011355.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.000) 5.31 ng/ul m

response 16435

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	105.61#
41.00	54.40	58.36
42.00	35.50	32.92

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	99968	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.375	136	423430	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	244319	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	492827	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	461236	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	463214	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	16072	4.929	ng/uL	0.00
4) Pyridine-d5	3.517	84	17202	2.091	ng/ul	0.02
7) Phenol-d5	6.769	99	70326	7.949	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	196098	31.261	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	189140	27.248	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	124634	17.942	ng/ul	-0.01
21) Nitrobenzene-d5	8.746	128	112741	39.122	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.469	143	108811	41.242	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	174931	32.832	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	78980	8.479	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	682409	40.255	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	765614	37.519	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	16435m	5.308	ng/ul	0.00
60) Fluorene-d10	15.263	176	564650	38.948	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	93063	40.655	ng/ul	0.00
73) Anthracene-d10	17.128	188	958130	43.253	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1104222	43.454	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1045879	44.772	ng/ul	0.00

Target Compounds Qvalue

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

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