

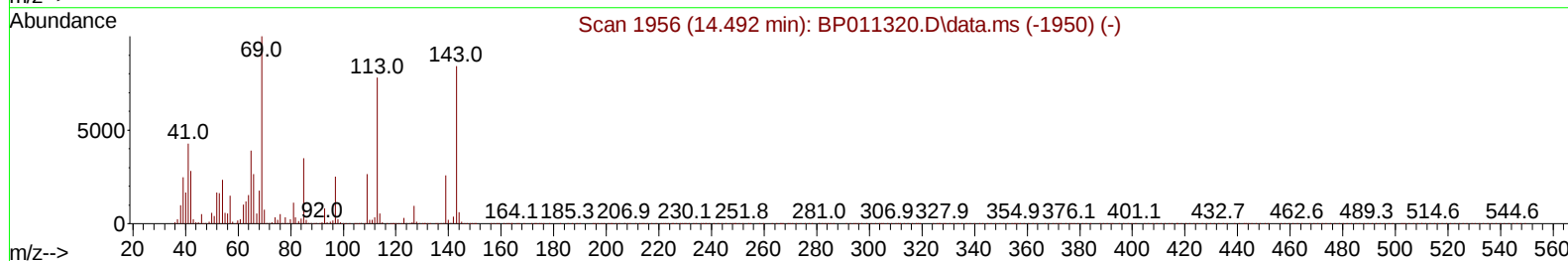
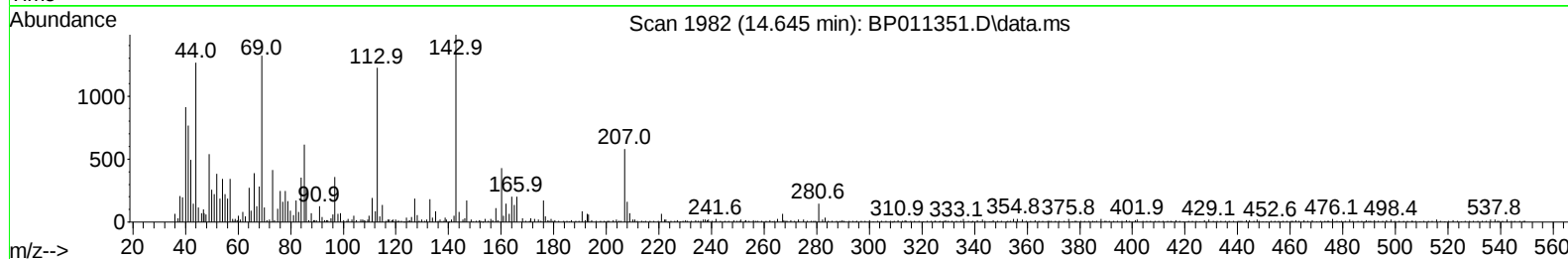
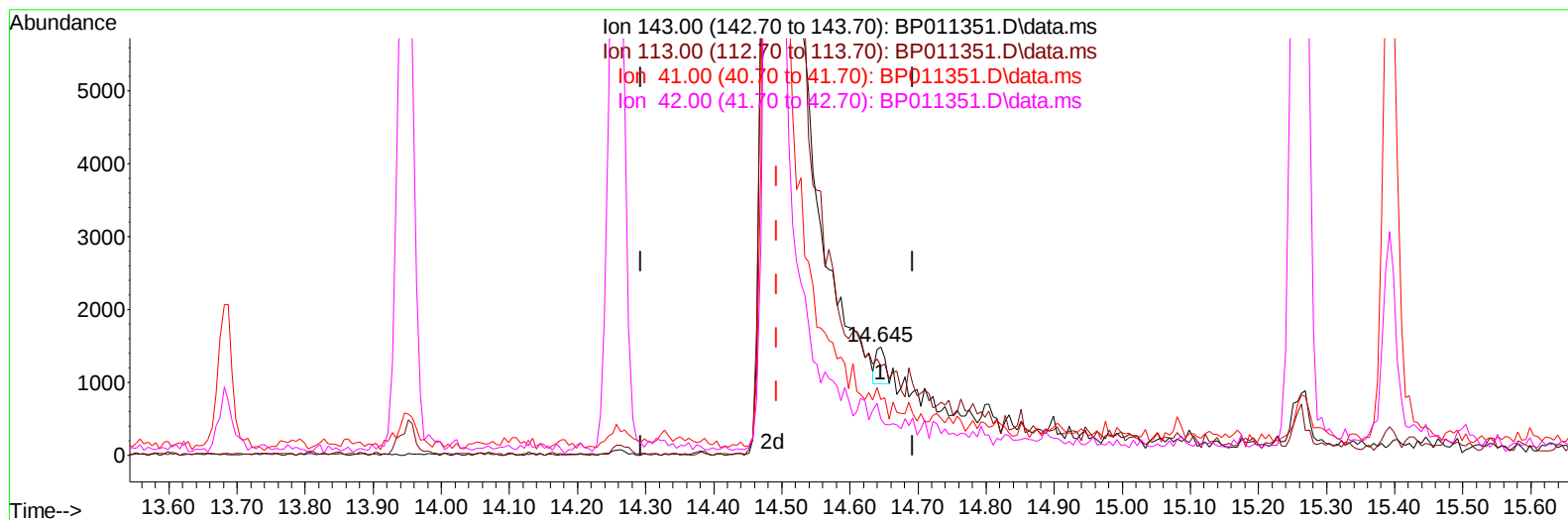
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011351.D
Acq On : 10 Aug 2022 04:16
Operator : CG/JU
Sample : PB146834BL
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK834

Manual IntegrationsAPPROVED

Quant Time: Aug 10 05:00:51 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: BP011351.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.153) 0.12 ng/u1

response 472

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	82.31
41.00	54.40	51.51
42.00	35.50	33.15

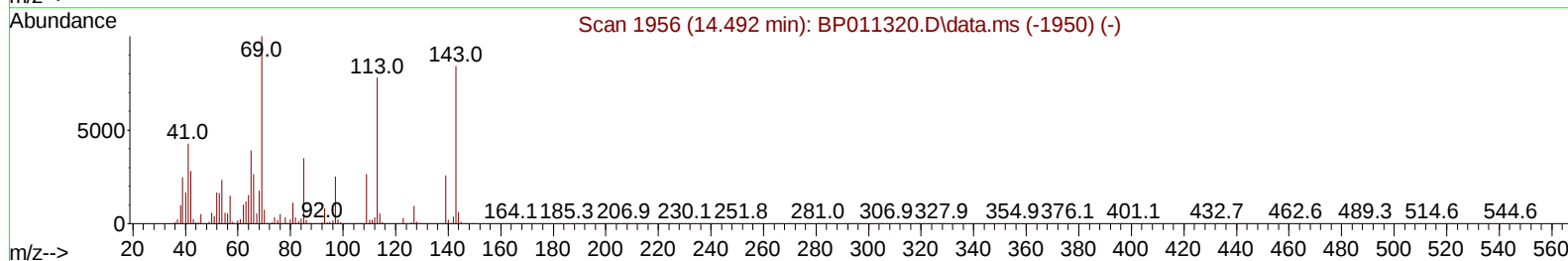
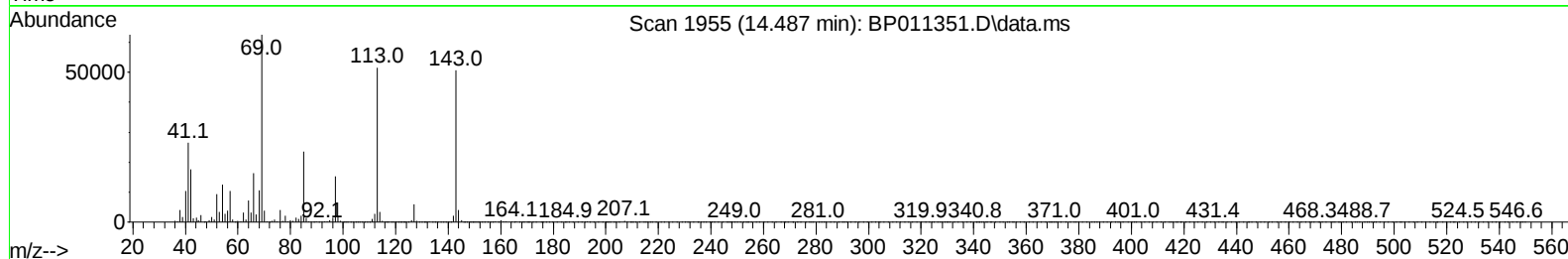
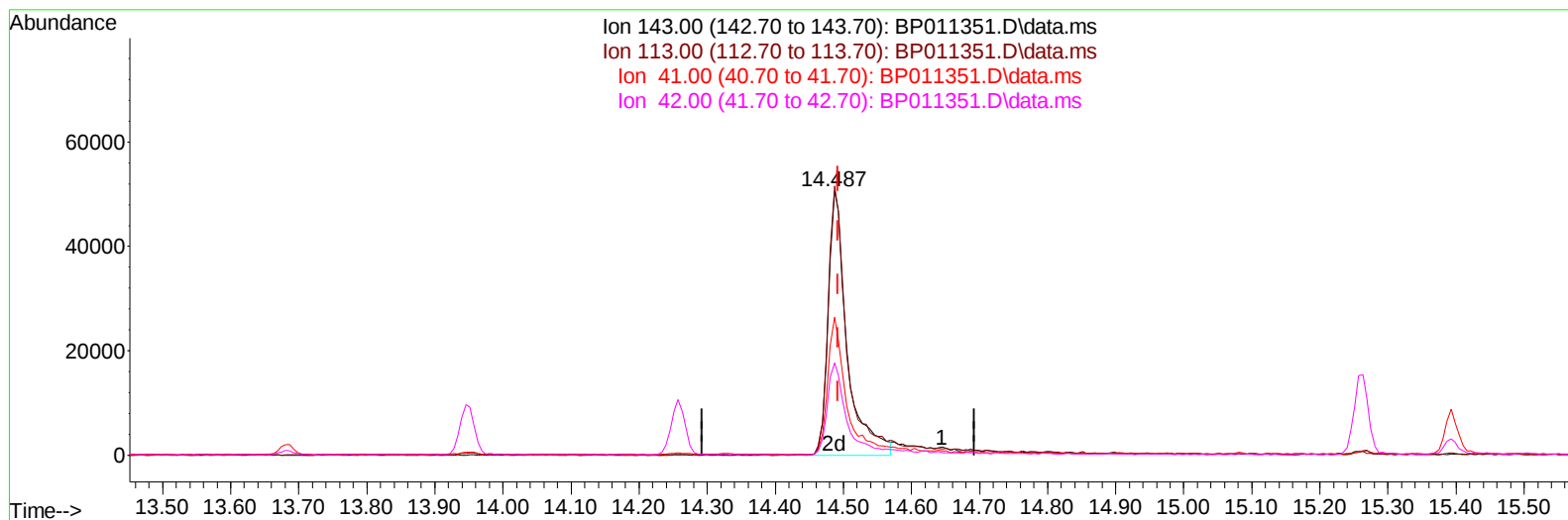
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TIC: BP011351.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (-0.006) 24.54 ng/ul m

response 97474

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	101.93#
41.00	54.40	52.21
42.00	35.50	34.94

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	131622	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	553398	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	313443	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	626775	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	579930	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	585592	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	17007	3.961	ng/uL	0.00
4) Pyridine-d5	3.505	84	237002	21.883	ng/ul	0.00
7) Phenol-d5	6.769	99	276642	23.750	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	186647	22.599	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	223728	24.479	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	226700	24.787	ng/ul	-0.01
21) Nitrobenzene-d5	8.752	128	106957	28.399	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	110313	31.991	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	183674	26.377	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	283542	23.291	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	704421	32.390	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	773844	29.559	ng/ul	0.00
54) 4-Nitrophenol-d4	14.487	143	97474m	24.539	ng/ul	0.00
60) Fluorene-d10	15.263	176	545504	29.330	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	86718	29.787	ng/ul	0.00
73) Anthracene-d10	17.128	188	898446	31.891	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1040886	32.578	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	974359	32.994	ng/ul	-0.01

Target Compounds Qvalue

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

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