

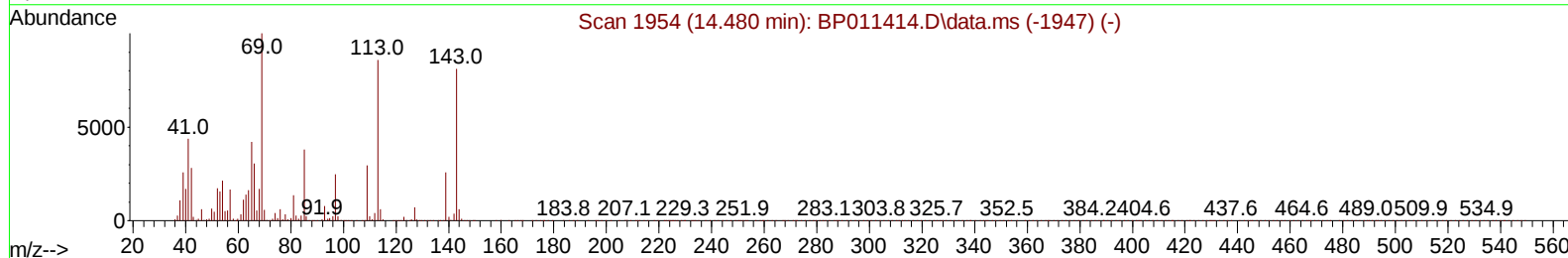
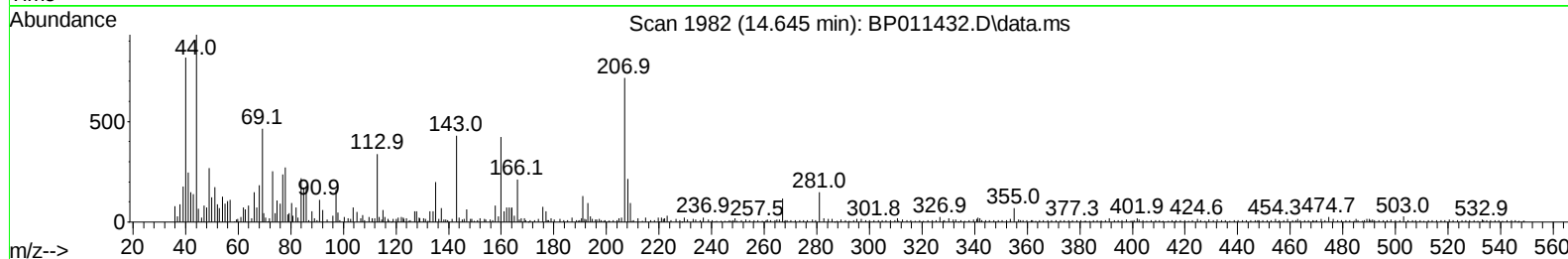
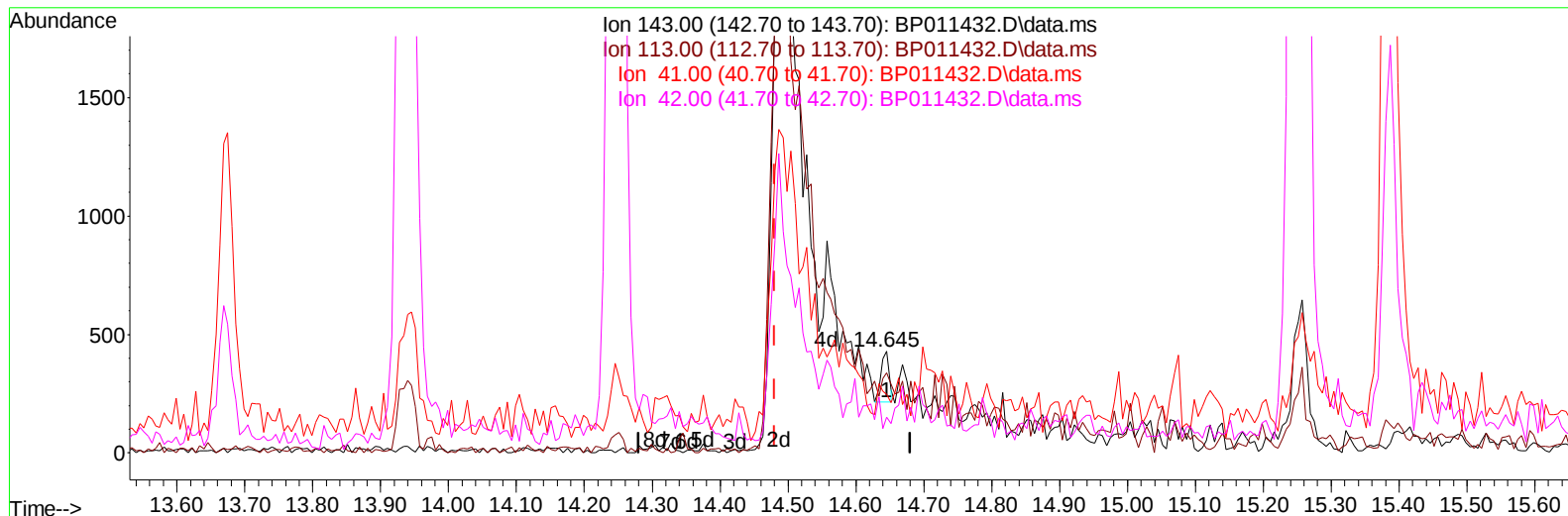
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011432.D
Acq On : 12 Aug 2022 19:50
Operator : CG/JU
Sample : N4165-15
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGBZ3

Manual IntegrationsAPPROVED

Quant Time: Aug 12 23:17:42 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 :Yogesh Patel 08/13/2022
Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011432.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.165) 0.08 ng/u1

response 179

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	84.42
41.00	54.40	57.44
42.00	35.50	34.88

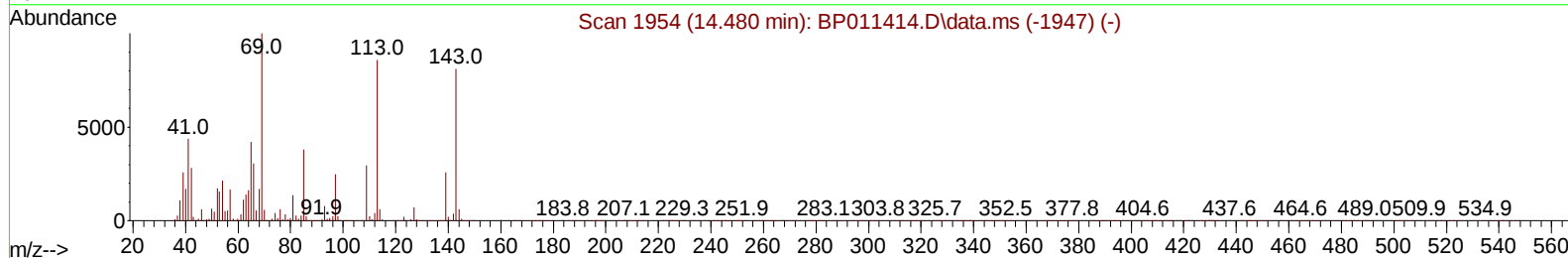
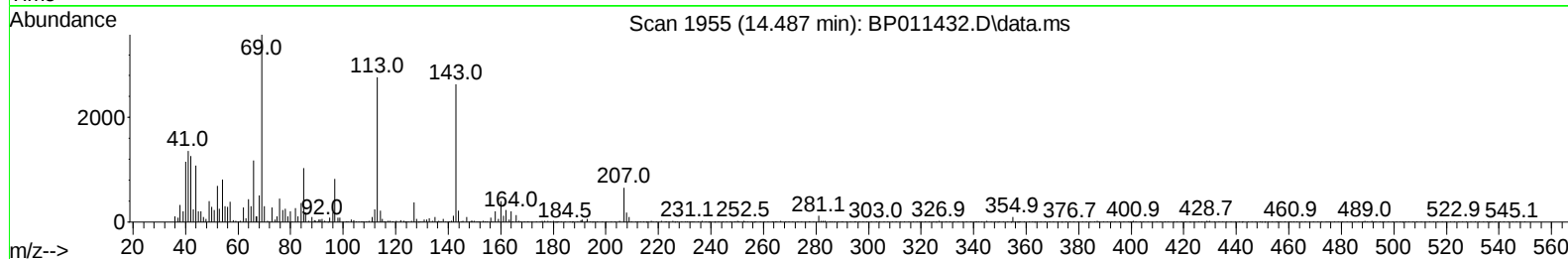
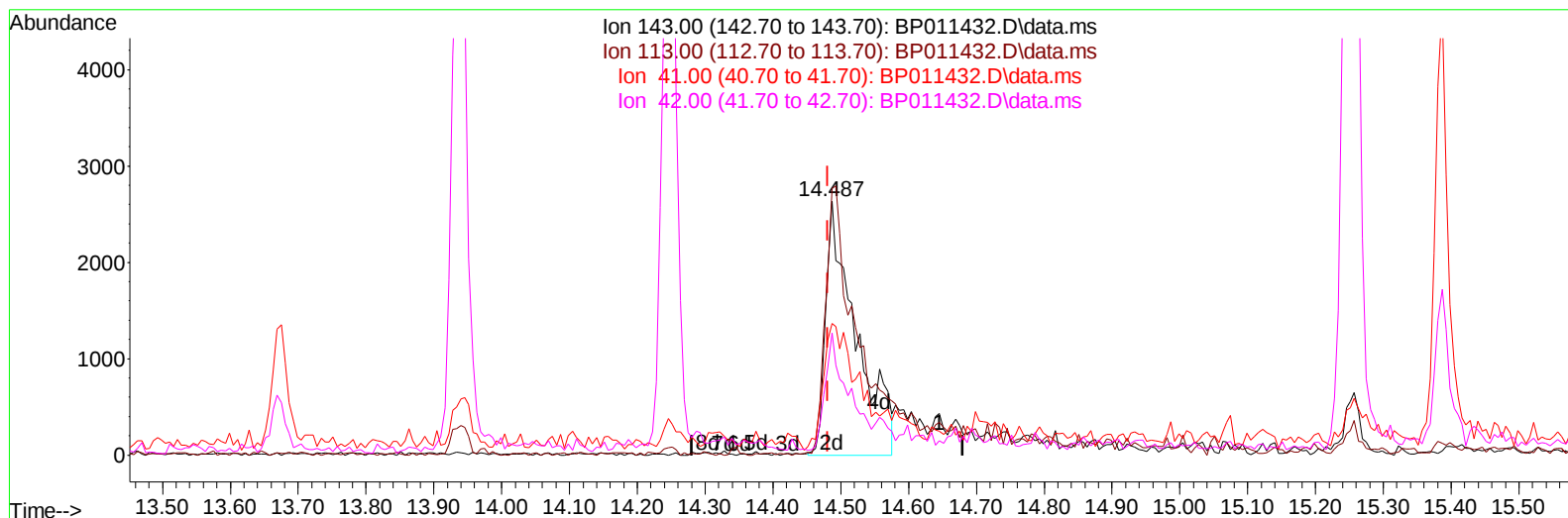
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(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.006) 3.50 ng/ul m

response 8139

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	104.97#
41.00	54.40	51.80
42.00	35.50	47.93#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	75995	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	319291	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.246	164	183557	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	379120	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	374555	20.000	ng/ul	# 0.00
88) Perylene-d12	23.457	264	385227	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	10817	4.364	ng/uL	0.00
4) Pyridine-d5	3.493	84	51745	8.275	ng/ul	0.00
7) Phenol-d5	6.758	99	46101	6.855	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	131596	27.596	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	123340	23.374	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	77727	14.719	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	71486	32.897	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	66728	33.540	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	103966	25.877	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	144447	20.565	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	410575	32.237	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	443290	28.914	ng/ul	0.00
54) 4-Nitrophenol-d4	14.487	143	8139m	3.499	ng/ul	0.00
60) Fluorene-d10	15.251	176	339993	31.215	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	47022	26.702	ng/ul	0.00
73) Anthracene-d10	17.122	188	598558	35.125	ng/ul	0.00
81) Pyrene-d10	19.386	212	675247	32.722	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	668053	34.387	ng/ul	0.00

Target Compounds Qvalue

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

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