

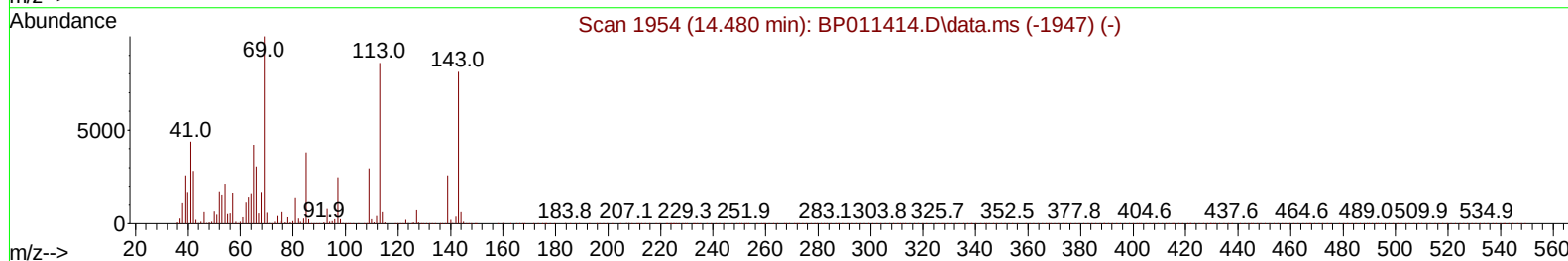
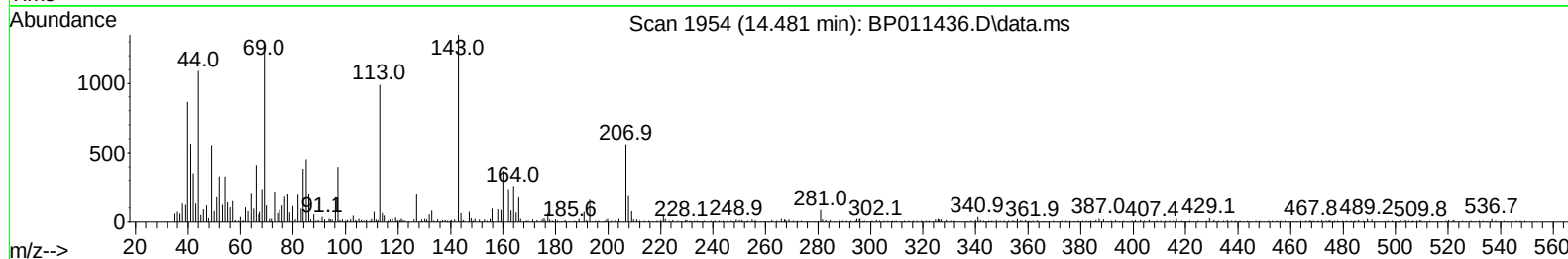
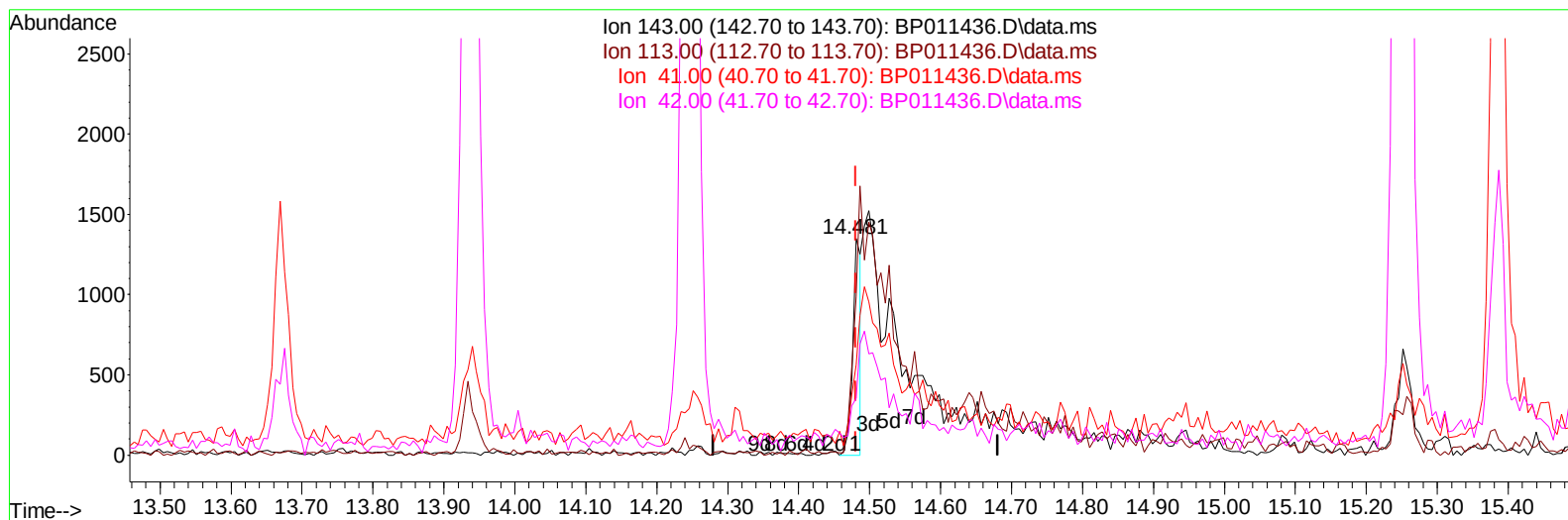
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011436.D
 Acq On : 12 Aug 2022 22:08
 Operator : CG/JU
 Sample : N4165-19
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGC13

Manual IntegrationsAPPROVED

Quant Time: Aug 12 23:19:13 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: BP011436.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.000) 0.43 ng/ul

response 1136

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	73.26
41.00	54.40	41.85#
42.00	35.50	26.07#

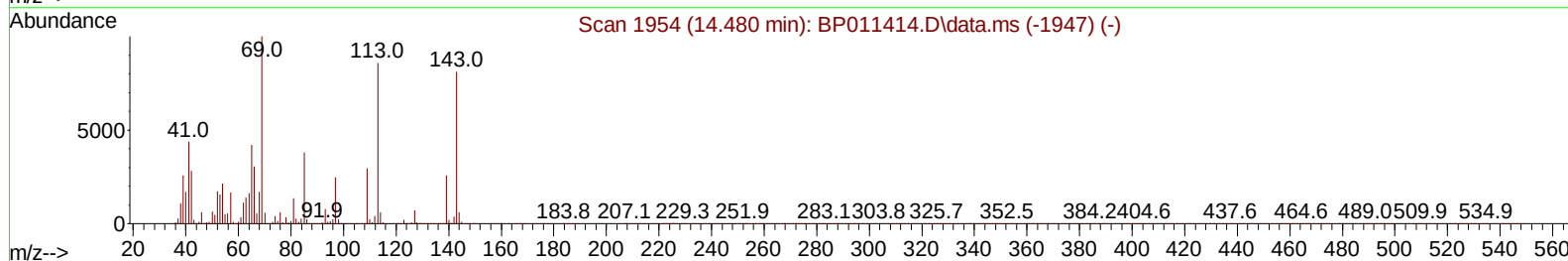
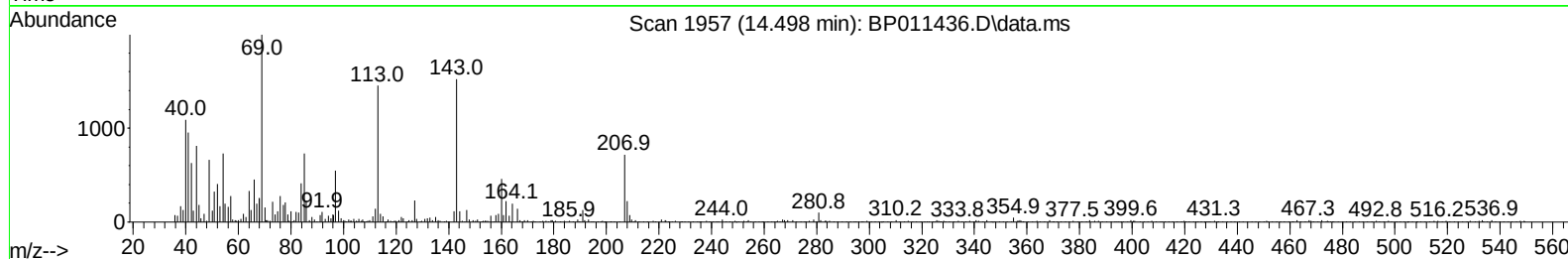
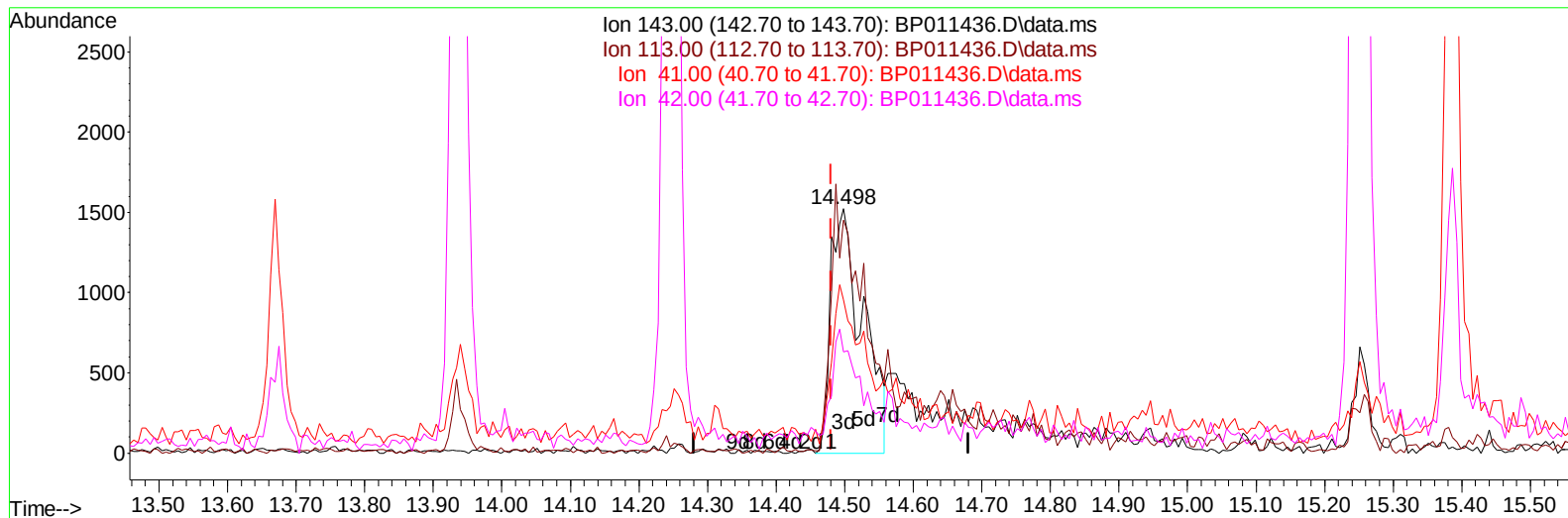
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TIC: BP011436.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 1.85 ng/ul m

response 4950

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	95.21#
41.00	54.40	62.69
42.00	35.50	41.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	81971	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	360566	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.246	164	210728	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	435903	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	428568	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	453643	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	10118	3.784	ng/uL	0.00
4) Pyridine-d5	3.499	84	36843	5.462	ng/ul	0.01
7) Phenol-d5	6.758	99	33761	4.654	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	131191	25.506	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	111507	19.591	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	64997	11.411	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	73388	29.907	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	67852	30.201	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	105152	23.176	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	144012	18.156	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	439501	30.059	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	476105	27.051	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	4950m	1.854	ng/ul	0.02
60) Fluorene-d10	15.251	176	358350	28.658	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	51557	25.464	ng/ul	0.00
73) Anthracene-d10	17.122	188	659565	33.663	ng/ul	0.00
81) Pyrene-d10	19.386	212	753770	31.924	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	749614	32.766	ng/ul	0.00

Target Compounds Qvalue

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

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