

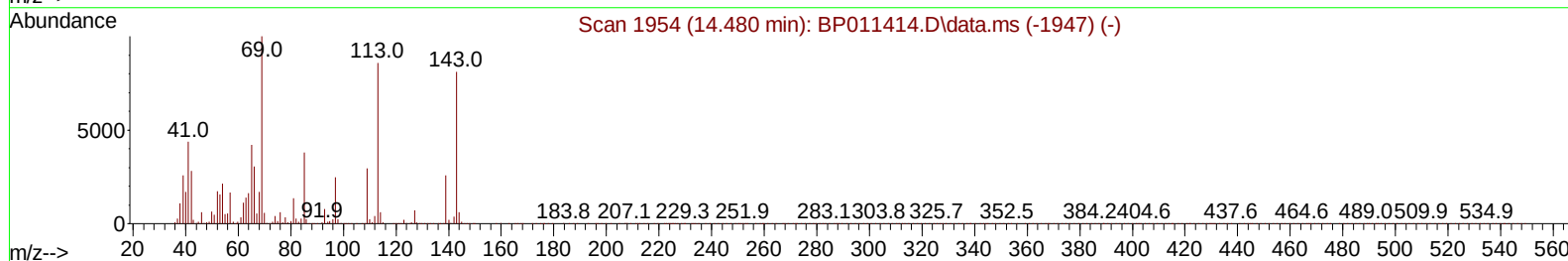
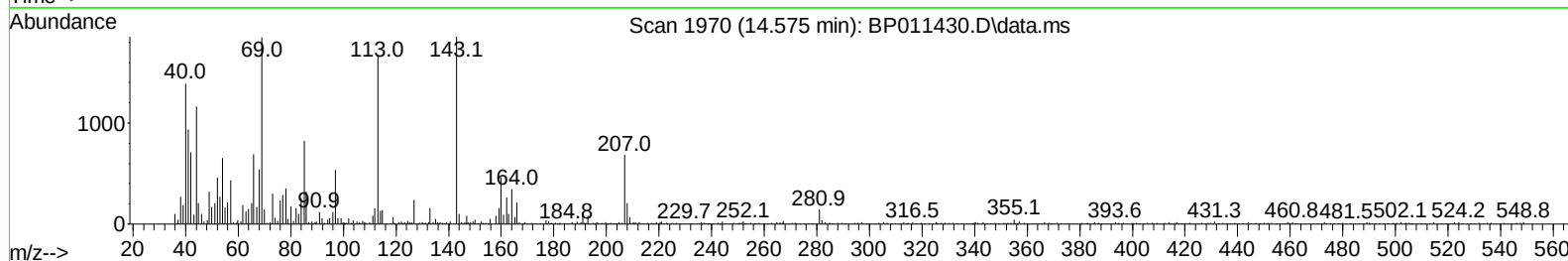
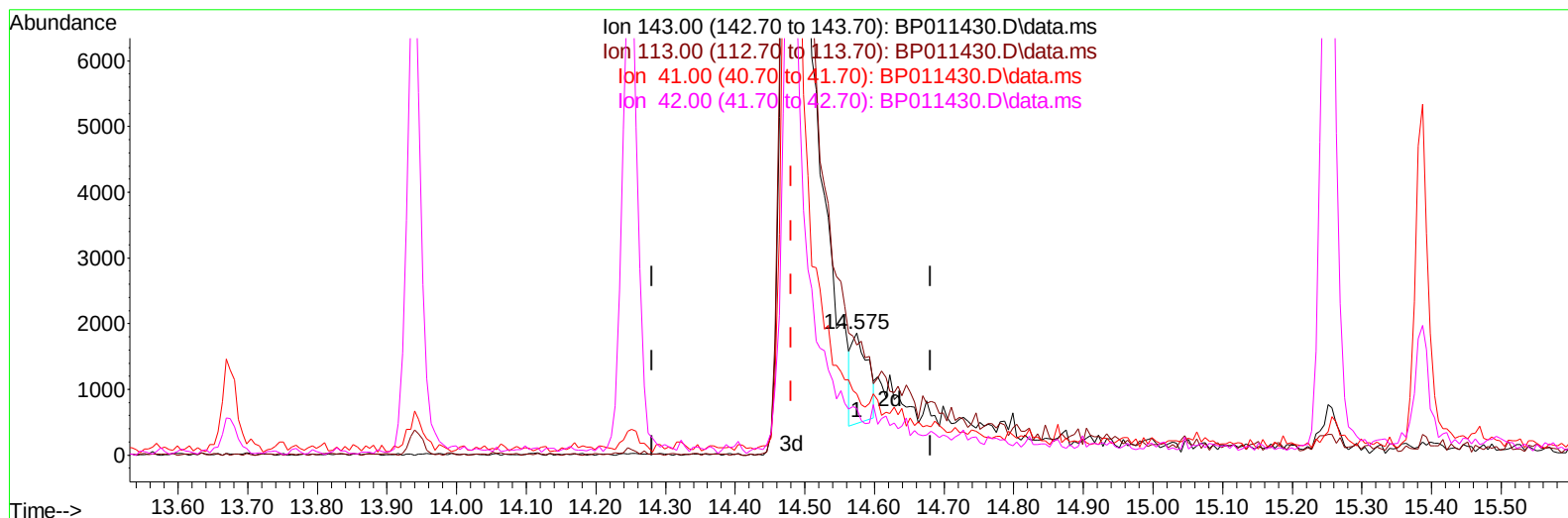
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011430.D
Acq On : 12 Aug 2022 18:41
Operator : CG/JU
Sample : PB146960BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK960

Manual IntegrationsAPPROVED

Quant Time: Aug 12 23:16:55 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: BP011430.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.094) 0.82 ng/ul

response 2180

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	90.13
41.00	54.40	50.46
42.00	35.50	38.38

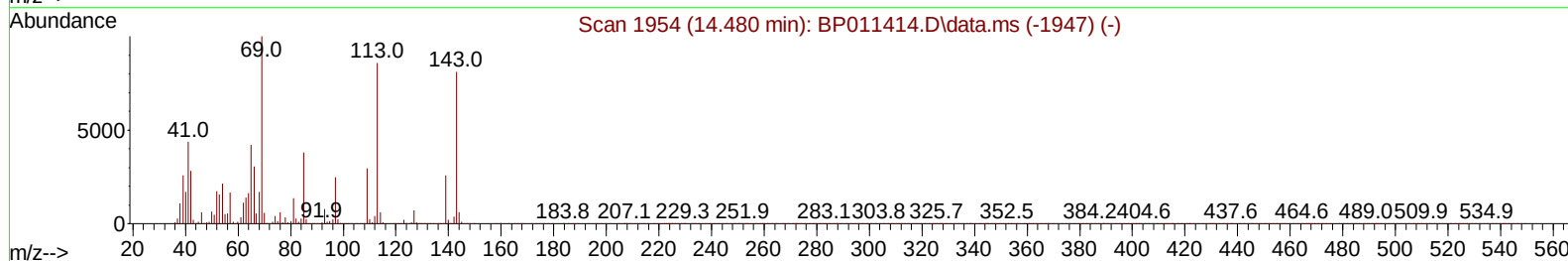
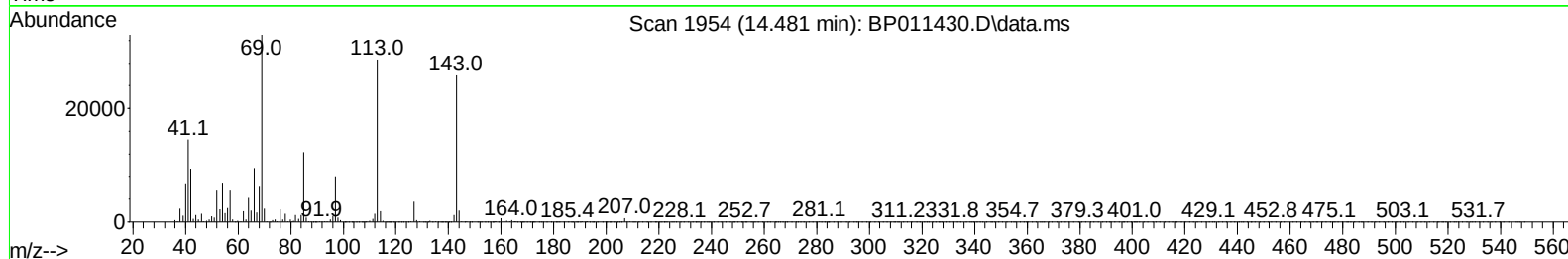
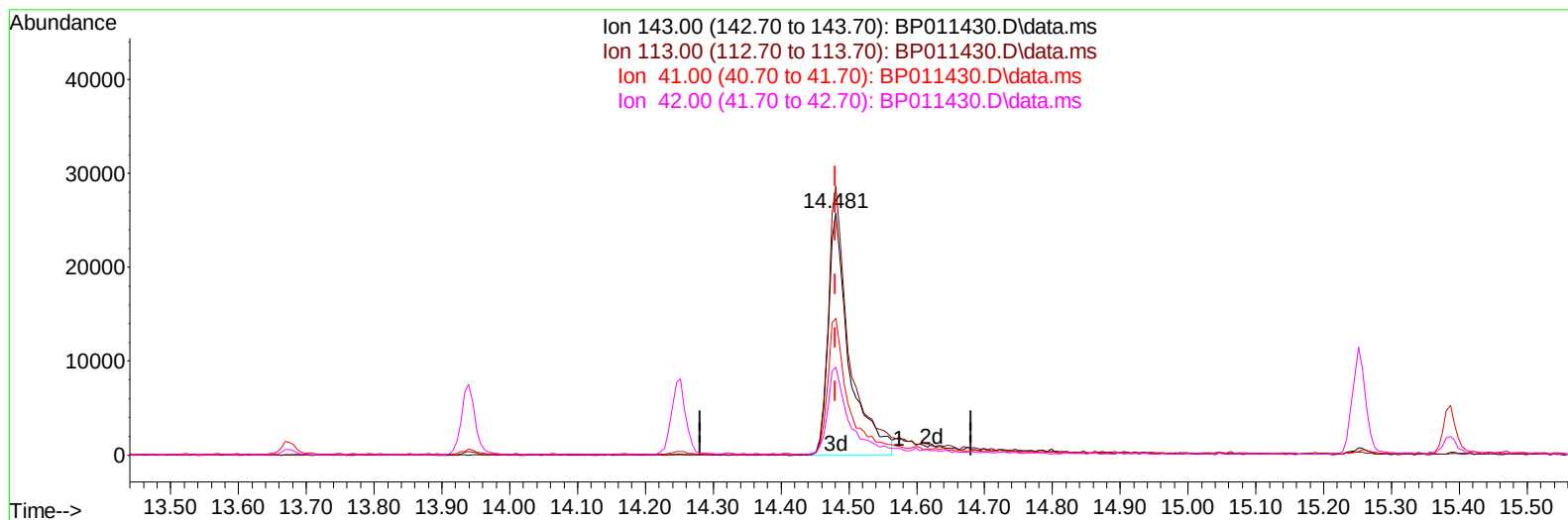
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TIC: BP011430.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.000) 20.89 ng/ul m

response 55374

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	111.04#
41.00	54.40	56.33
42.00	35.50	36.39

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	86210	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	366026	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.251	164	209192	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	436268	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.151	240	418926	20.000	ng/ul	0.00
88) Perylene-d12	23.463	264	450825	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	14369	5.110	ng/uL	0.00
4) Pyridine-d5	3.493	84	92310	13.013	ng/ul	0.00
7) Phenol-d5	6.758	99	196388	25.741	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	138817	25.661	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	163358	27.289	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	152275	25.420	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	76924	30.880	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	75863	33.263	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	118265	25.678	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	102353	12.711	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	432770	29.816	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	478315	27.376	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	55374m	20.887	ng/ul	0.00
60) Fluorene-d10	15.251	176	353795	28.502	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	52259	25.789	ng/ul	0.00
73) Anthracene-d10	17.122	188	596886	30.439	ng/ul	0.00
81) Pyrene-d10	19.386	212	671367	29.089	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	660456	29.050	ng/ul	0.00

Target Compounds Qvalue

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

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