

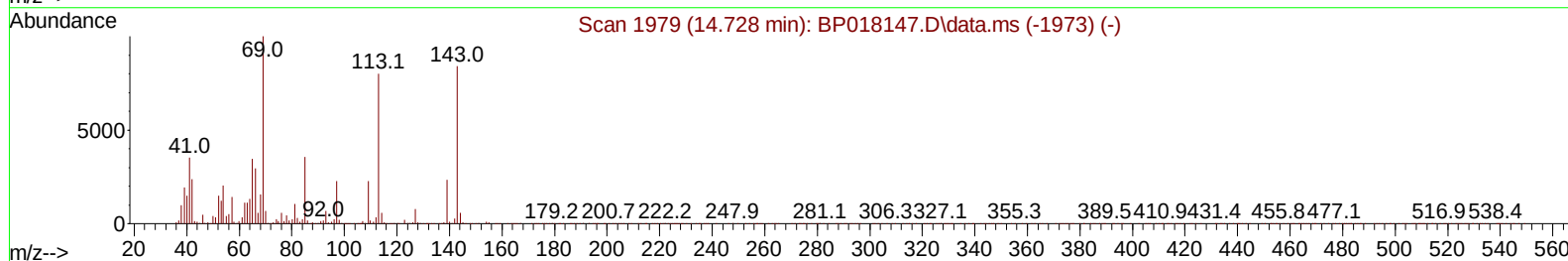
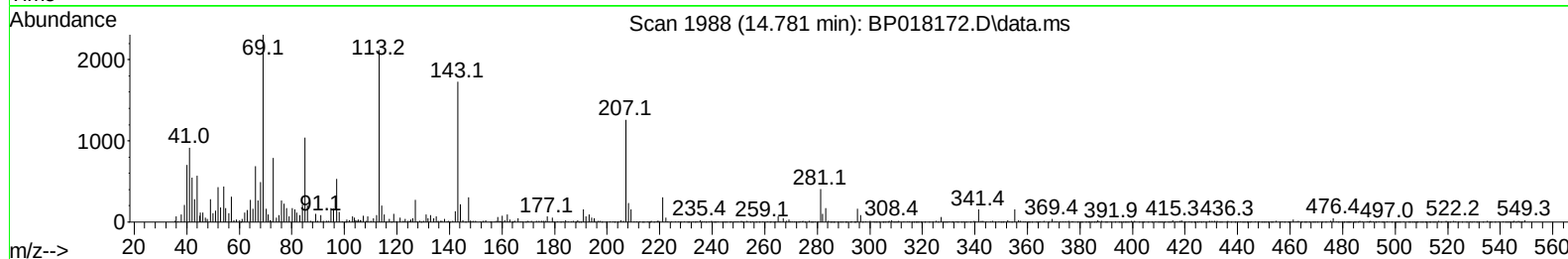
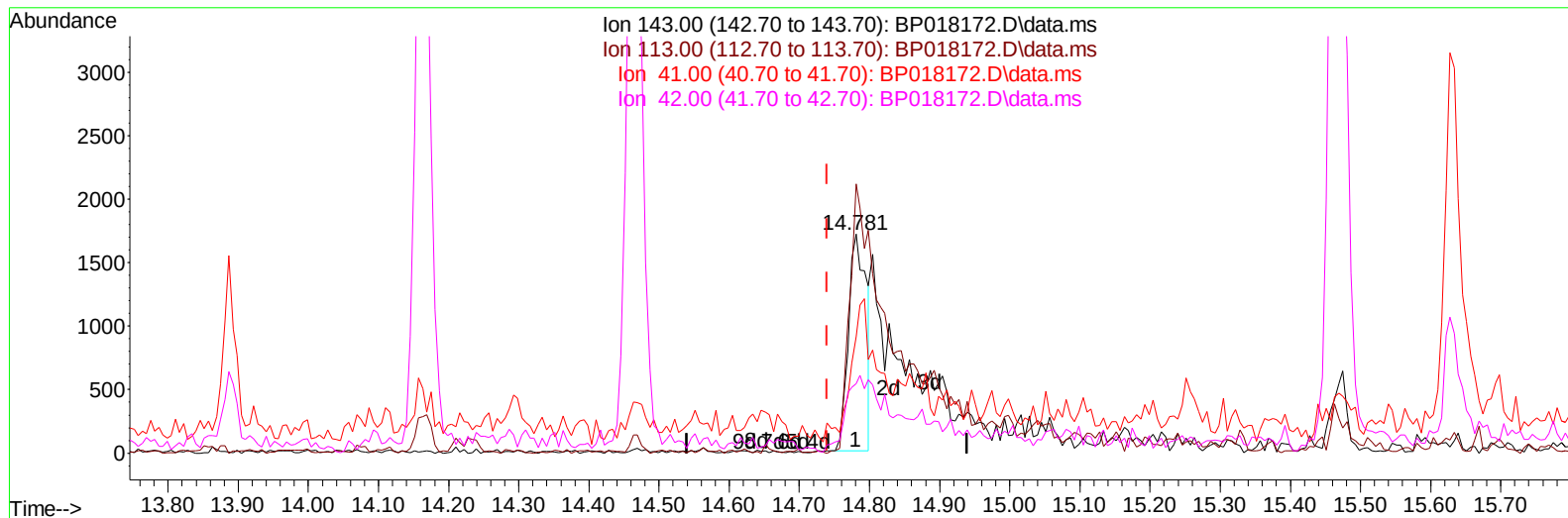
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018172.D
Acq On : 15 Nov 2023 00:37
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH362

Manual IntegrationsAPPROVED

Quant Time: Nov 15 01:54:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023



TIC: BP018172.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.781min (+ 0.041) 1.35 ng/u1

response 3052

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	122.84
41.00	47.70	53.10
42.00	30.10	31.77

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	65860	20.000	ng/ul	0.00
20) Naphthalene-d8	10.605	136	263264	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	163375	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.245	188	380282	20.000	ng/ul	-0.01
79) Chrysene-d12	21.375	240	392023	20.000	ng/ul	-0.02
88) Perylene-d12	23.845	264	487537	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6260	3.353	ng/uL	0.00
4) Pyridine-d5	3.617	84	26908	5.528	ng/ul	0.01
7) Phenol-d5	6.963	99	32033	5.038	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	95681	25.606	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	94596	18.787	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	60320	11.577	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	57921	24.420	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	61487	22.882	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	101213	21.133	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	126799	19.306	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	431509	28.626	ng/ul	-0.01
49) Acenaphthylene-d8	14.163	160	416119	25.344	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	8211m	3.645	ng/ul	0.04
60) Fluorene-d10	15.469	176	355043	28.528	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	49851	18.133	ng/ul	-0.04
73) Anthracene-d10	17.345	188	589017	28.639	ng/ul	-0.01
81) Pyrene-d10	19.604	212	801491	30.608	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.680	264	892114	31.469	ng/ul	-0.02

Target Compounds Qvalue

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

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