

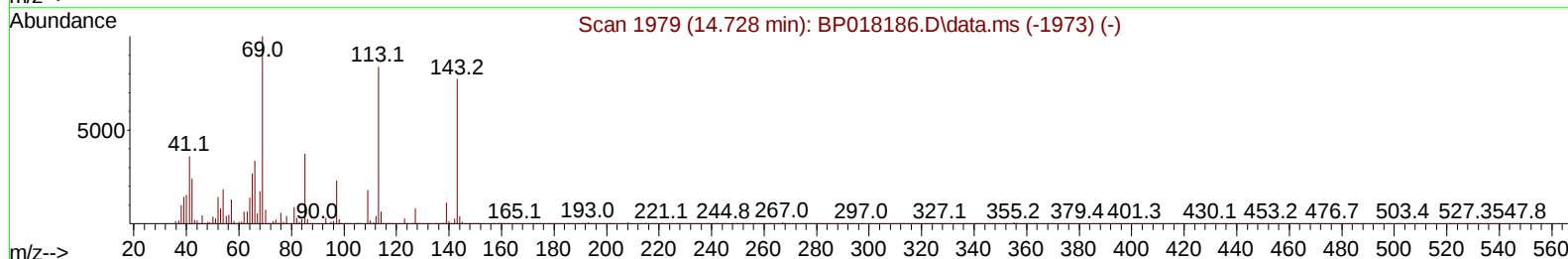
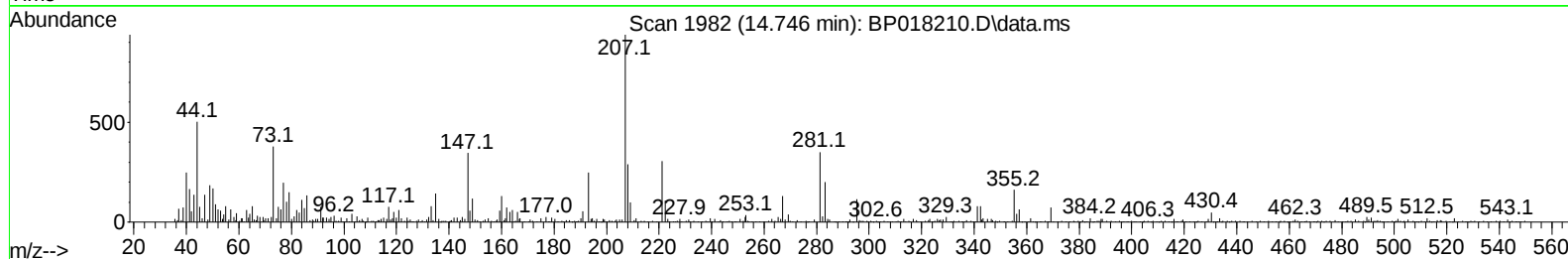
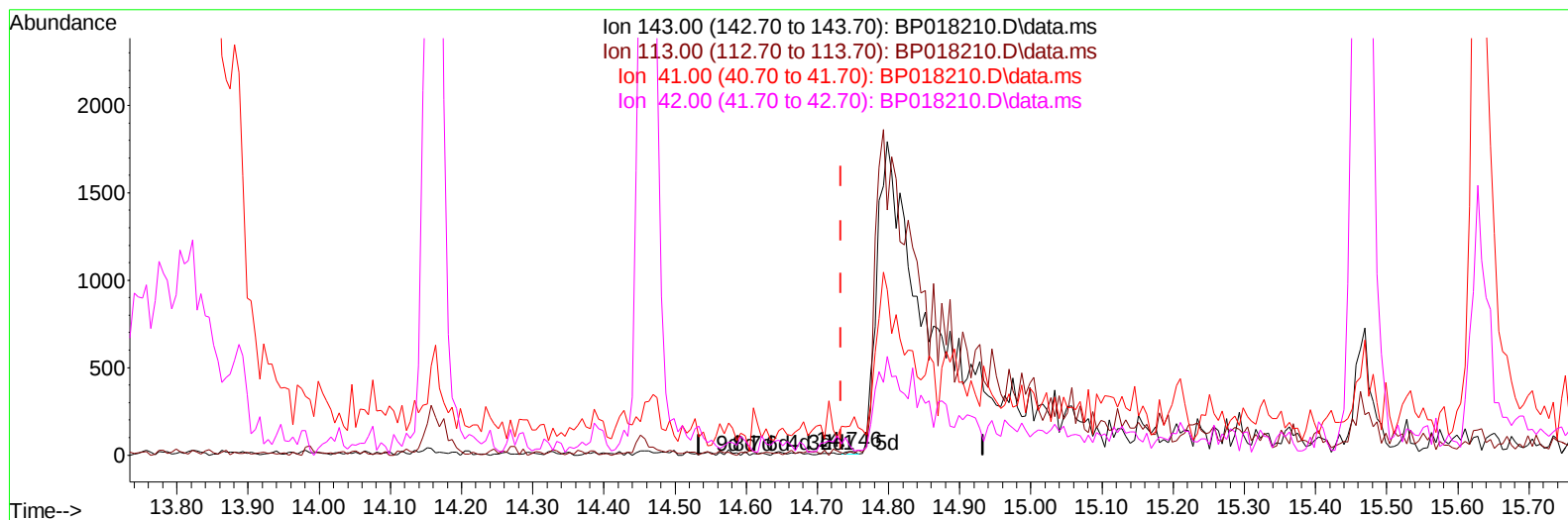
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDM0

Manual IntegrationsAPPROVED

Quant Time: Nov 16 21:38:57 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 18:13:37 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/17/2023
Supervised By :mohammad ahmed 11/17/2023



TIC: BP018210.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.746min (+ 0.012) 0.01 ng/u1

response 16

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	95.24
41.00	47.70	785.71#
42.00	30.10	257.14#

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Reviewed By :Yogesh Patel 11/17/2023
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Quant Time: Nov 16 22:09:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	69827	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	301056	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	184322	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	435692	20.000	ng/ul	0.00
79) Chrysene-d12	21.381	240	436233	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	488258	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6788	3.429	ng/uL	0.00
4) Pyridine-d5	3.605	84	36645	7.100	ng/ul	0.01
7) Phenol-d5	6.964	99	29598	4.391	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	87925	22.194	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	93596	17.533	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	58314	10.556	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	54964	20.264	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	65415	21.288	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	107076	19.551	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	77255	10.286	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	442834	26.038	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	431026	23.269	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	8872m	3.491	ng/ul	0.06
60) Fluorene-d10	15.463	176	357808	25.483	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	63877	20.280	ng/ul	0.00
73) Anthracene-d10	17.345	188	629506	26.715	ng/ul	0.00
81) Pyrene-d10	19.604	212	849750	29.162	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	823724	29.014	ng/ul	0.00

Target Compounds Qvalue

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

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