

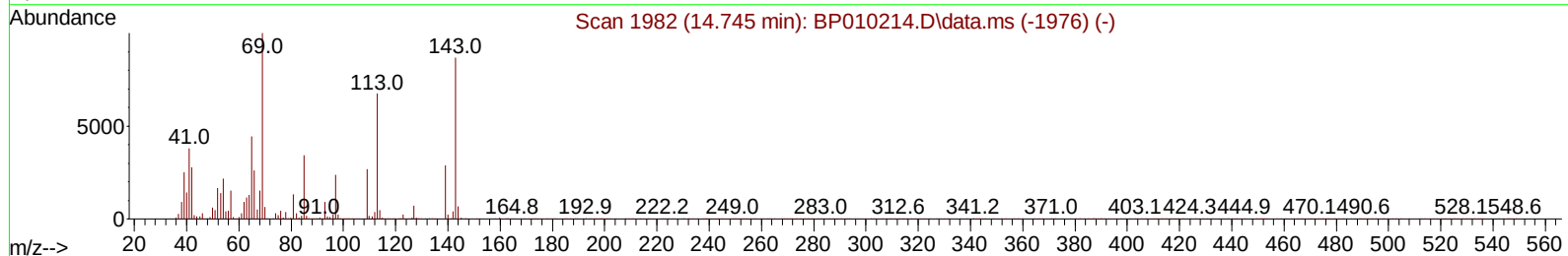
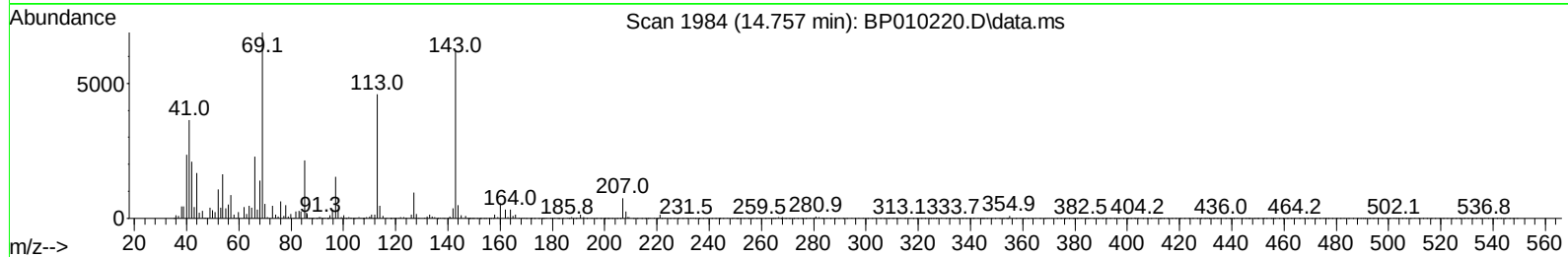
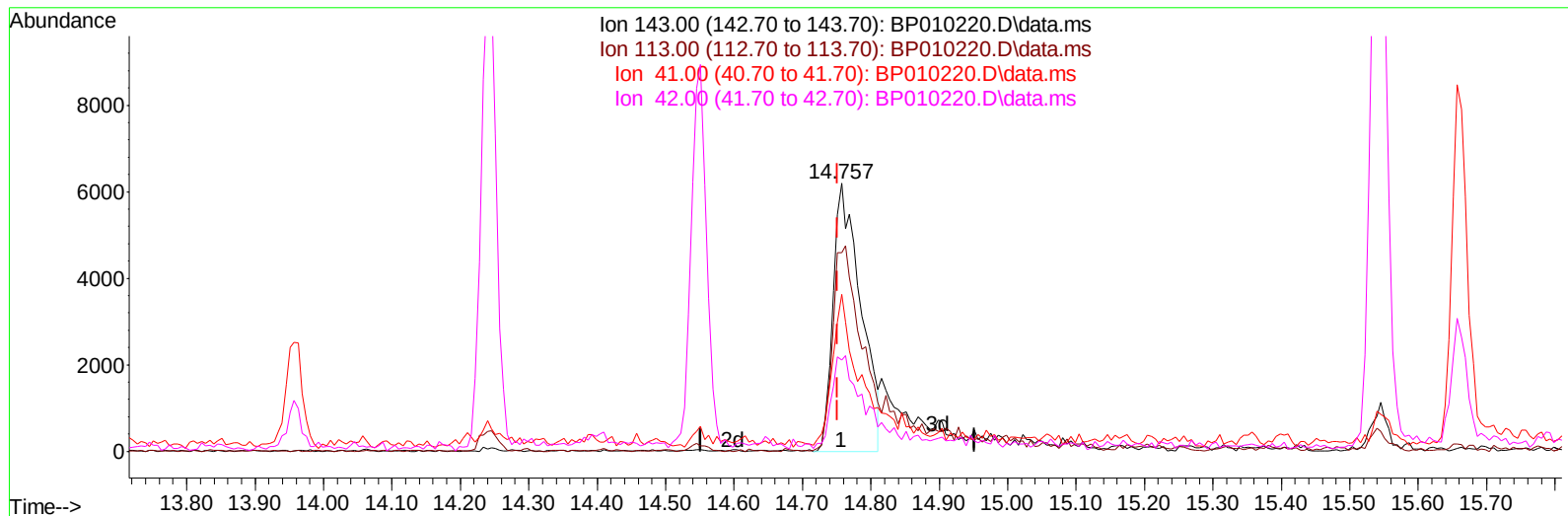
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010220.D
Acq On : 03 May 2022 22:53
Operator : CG/JU
Sample : N2670-03
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COW02

Manual IntegrationsAPPROVED

Quant Time: May 04 01:26:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/04/2022
Supervised By :Yogesh Patel 05/05/2022



TIC: BP010220.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 3.03 ng/ul

response 17365

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	74.21#
41.00	23.20	58.59#
42.00	18.00	34.03#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	155036	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	650787	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	399945	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	751168	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	668873	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	642397	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	18203	5.120	ng/uL	0.00
4) Pyridine-d5	3.781	84	73234	7.322	ng/ul	0.00
7) Phenol-d5	7.081	99	89086	6.317	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	314634	37.183	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	278216	26.639	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	176911	15.141	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	188146	36.350	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	192407	34.037	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	299360	27.203	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	392222	24.857	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1095075	35.199	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1259337	33.446	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.757	143	20690m	3.612	ng/ul	0.00
60) Fluorene-d10	15.545	176	946419	35.364	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	136389	28.715	ng/ul	0.00
73) Anthracene-d10	17.398	188	1431219	39.383	ng/ul	-0.01
81) Pyrene-d10	19.633	212	1591111	41.828	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1366045	40.562	ng/ul	-0.01

Target Compounds Qvalue

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

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