

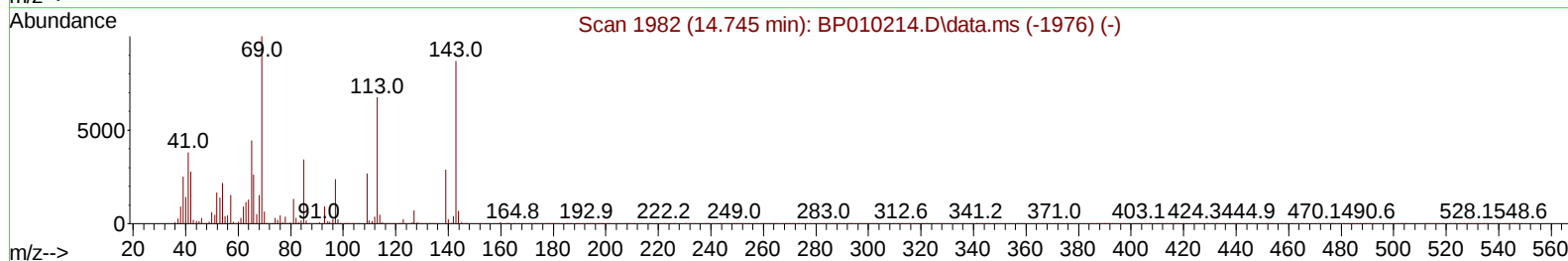
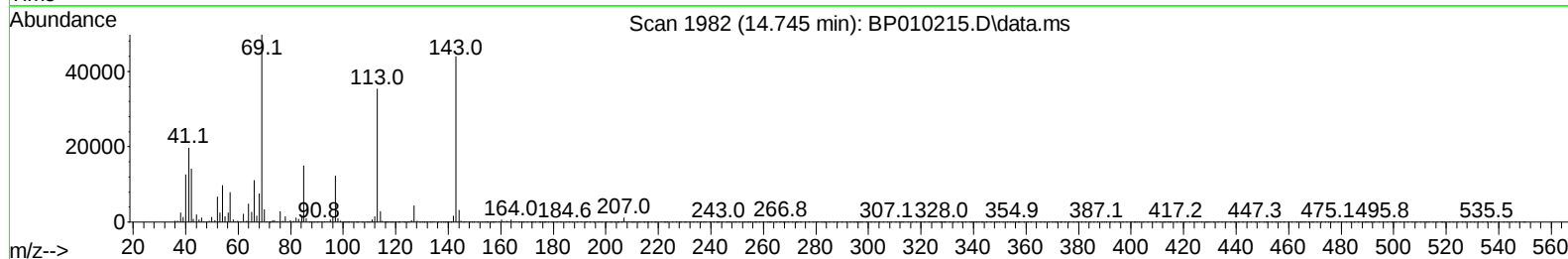
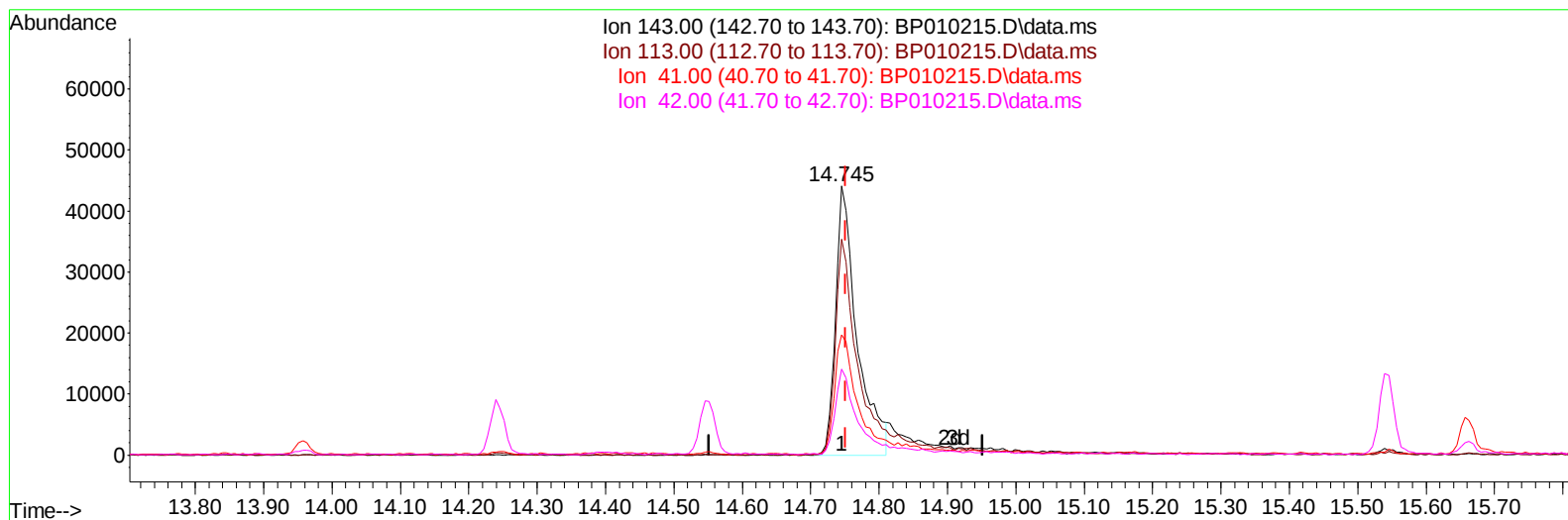
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010215.D
Acq On : 03 May 2022 20:05
Operator : CG/JU
Sample : PB144558BL
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK558

Manual IntegrationsAPPROVED

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

Quant Time: May 04 01:24:46 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



TIC: BP010215.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 16.16 ng/ul

response 96189

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.20#
41.00	23.20	44.77#
42.00	18.00	32.04#

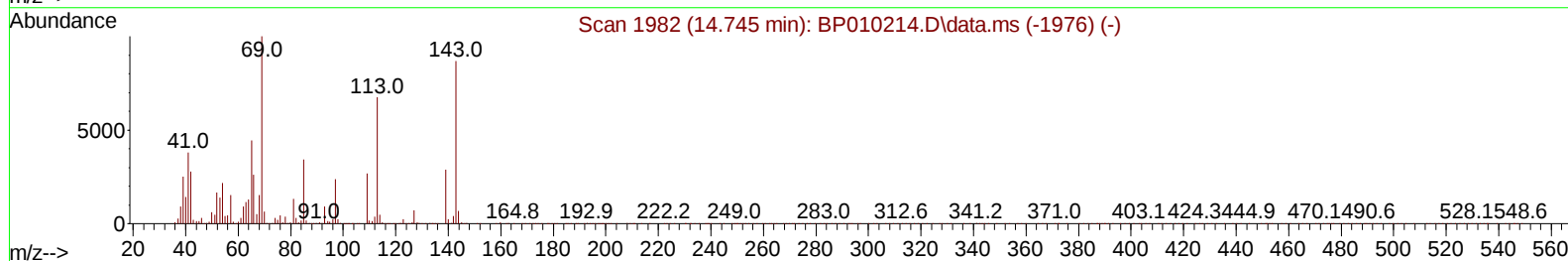
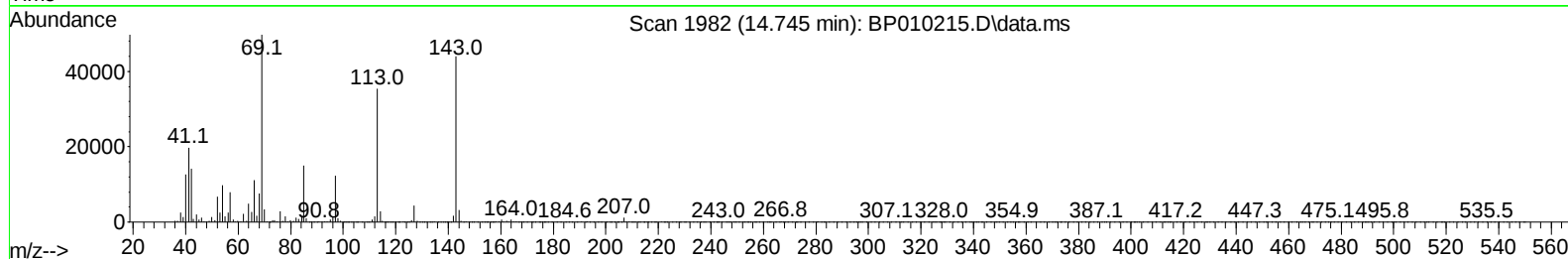
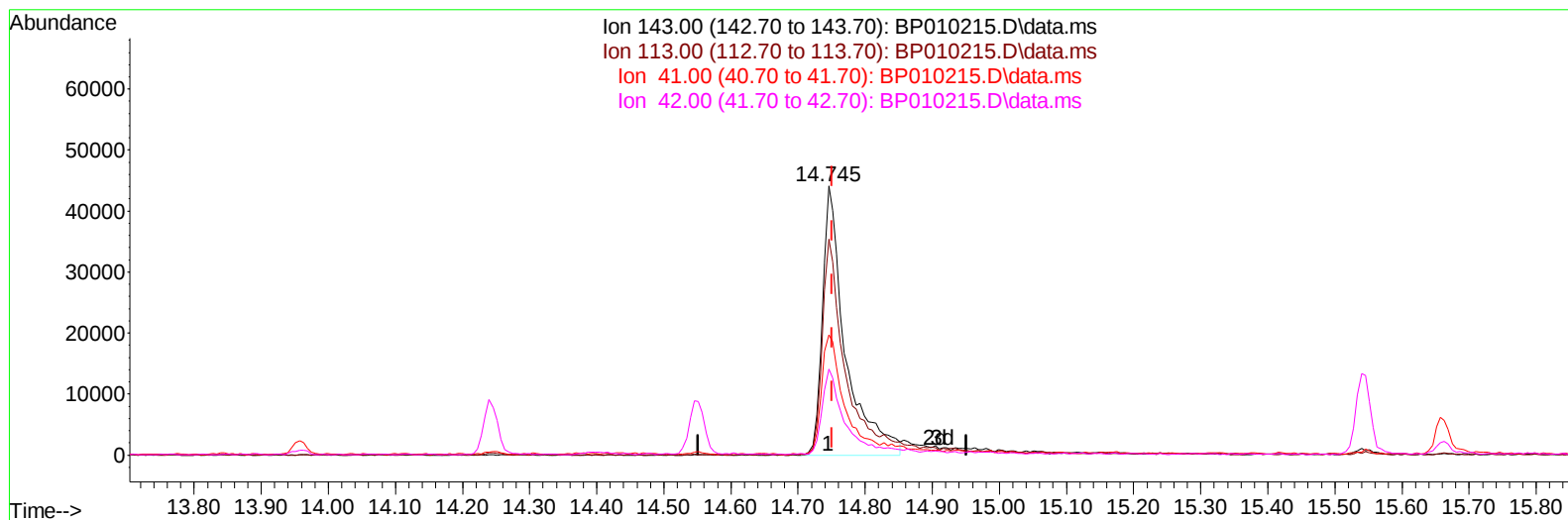
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TIC: BP010215.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 17.56 ng/ul m

response 104556

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.20#
41.00	23.20	44.77#
42.00	18.00	32.04#

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 Operator : CG/JU
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 Misc :
 ALS Vial : 52 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	163042	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.722	136	667557	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.551	164	415721	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	824280	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	728673	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	683108	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	20812	5.566	ng/uL	0.00
4) Pyridine-d5	3.770	84	277656	26.397	ng/ul	0.00
7) Phenol-d5	7.081	99	349900	23.594	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	251224	28.231	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	292299	26.614	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	289802	23.586	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	147376	27.758	ng/ul	0.00
24) 2-Nitrophenol-d4	9.805	143	156954	27.068	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	266457	23.605	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	377868	23.346	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	870582	26.921	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1003206	25.632	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	104556m	17.562	ng/ul	0.00
60) Fluorene-d10	15.545	176	743386	26.723	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	101036	19.385	ng/ul	0.00
73) Anthracene-d10	17.404	188	1137108	28.515	ng/ul	0.00
81) Pyrene-d10	19.633	212	1293070	31.204	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1063952	29.709	ng/ul	-0.01

Target Compounds Qvalue

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

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