

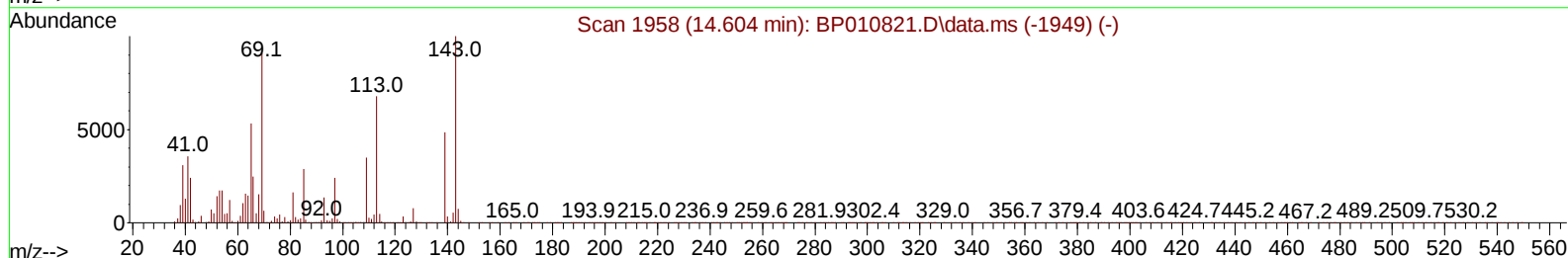
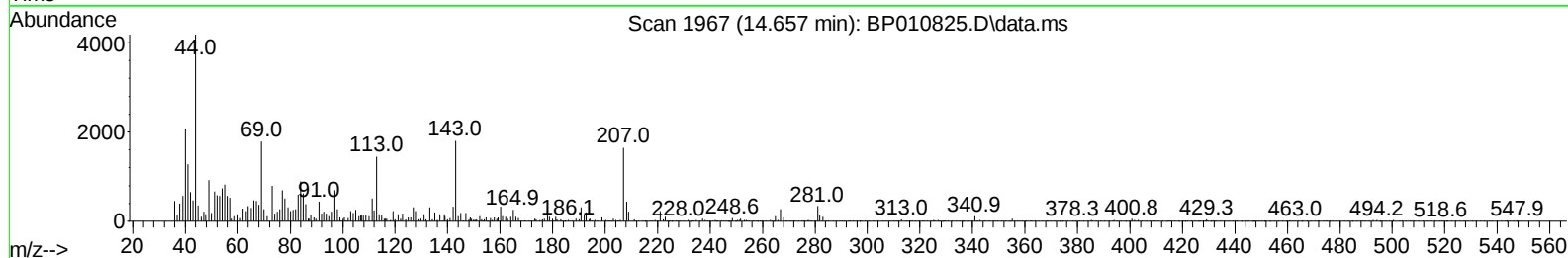
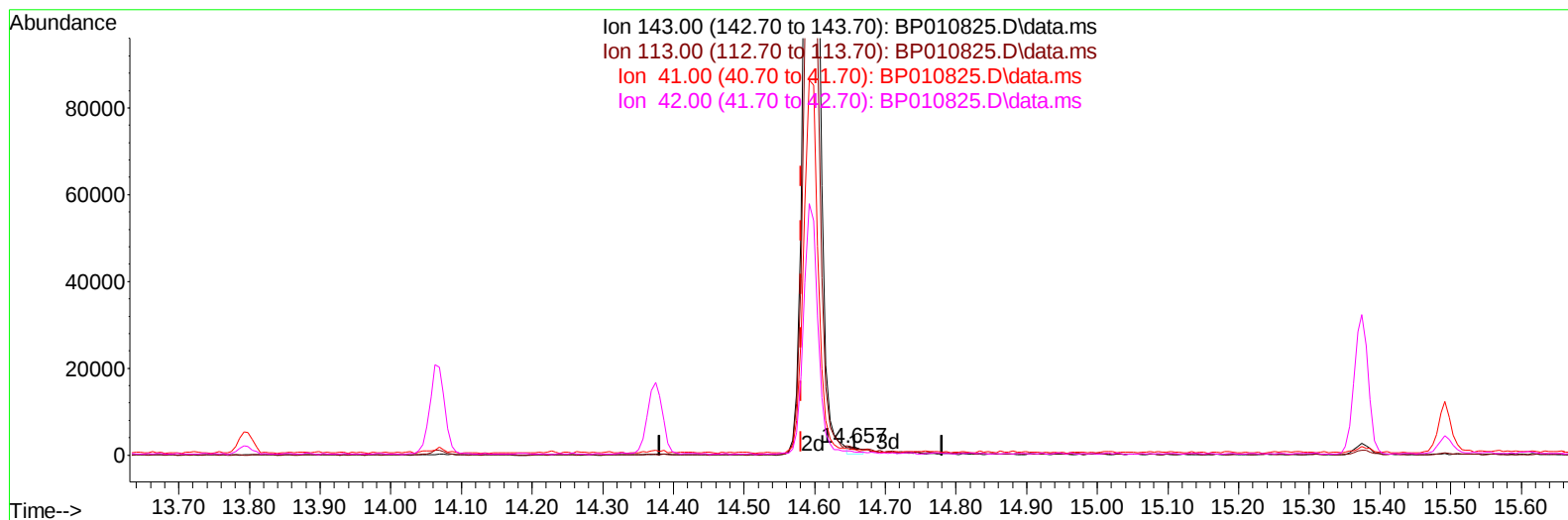
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010825.D
Acq On : 25 Jun 2022 13:21
Operator : CG/JU
Sample : PB145664BL
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK664

Manual IntegrationsAPPROVED

Quant Time: Jun 27 09:30:43 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 24 22:42:26 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/27/2022
Supervised By :mohammad ahmed 07/01/2022



TIC: BP010825.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.657min (+ 0.076) 0.21 ng/u1

response 1938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	80.23
41.00	59.00	70.60
42.00	41.10	36.43

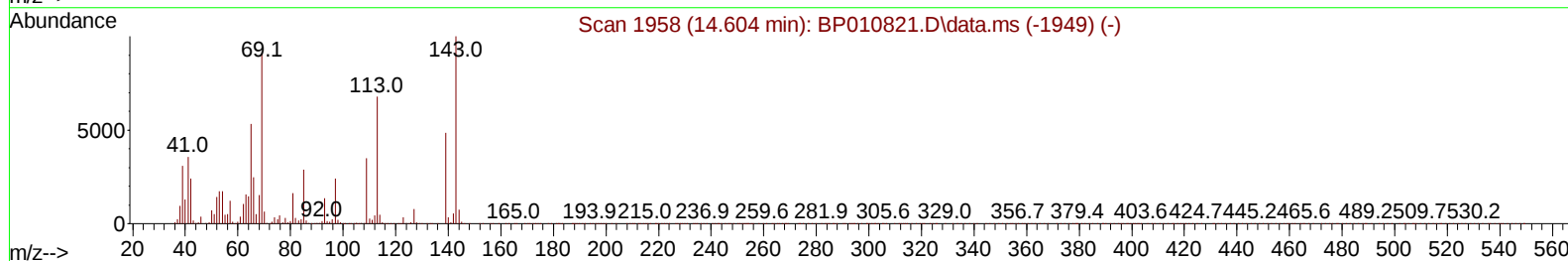
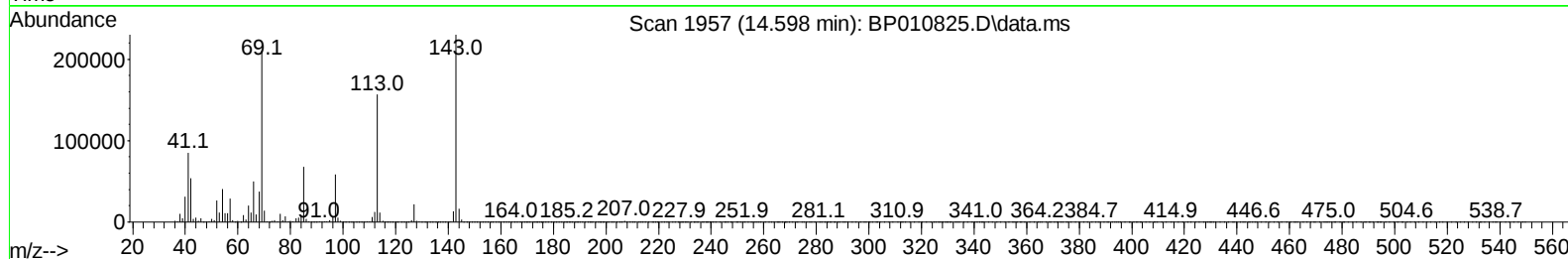
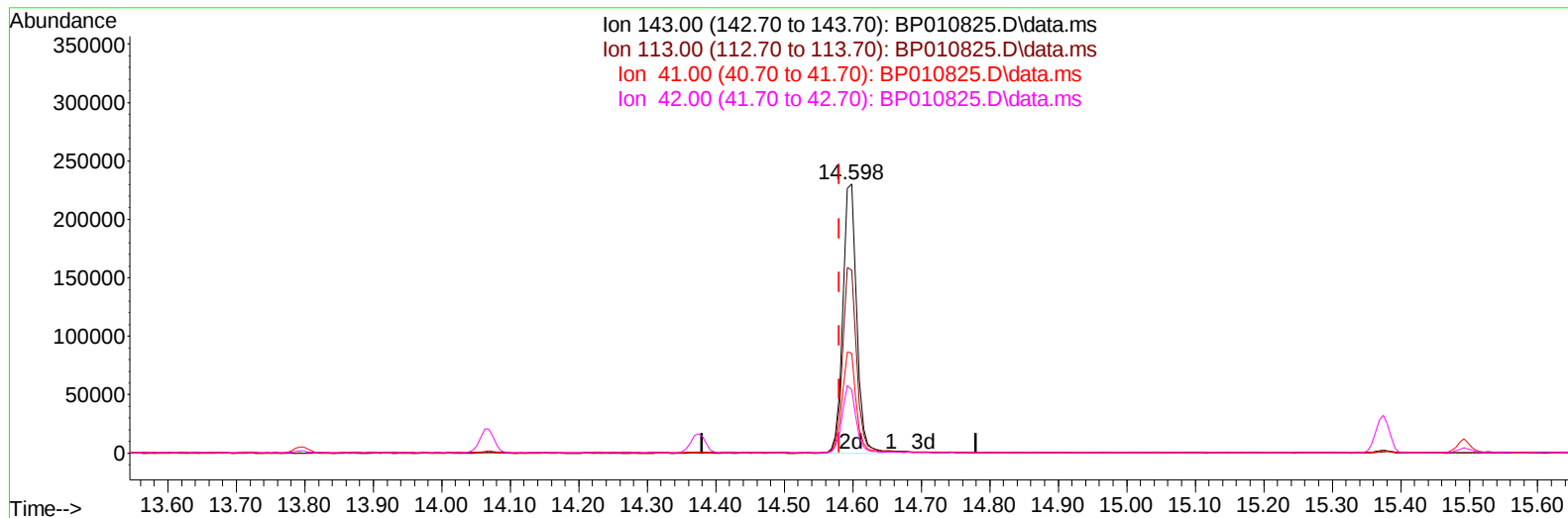
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TIC: BP010825.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.017) 34.94 ng/ul m

response 323132

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	68.10#
41.00	59.00	37.05#
42.00	41.10	23.49#

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 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	402186	20.000	ng/ul	0.00
20) Naphthalene-d8	10.516	136	1621113	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.375	164	864336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.139	188	1535499	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.251	240	1419877	20.000	ng/ul	# 0.00
88) Perylene-d12	23.621	264	1577719	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	77625	7.301	ng/uL	0.00
4) Pyridine-d5	3.616	84	1046087	36.963	ng/ul	0.01
7) Phenol-d5	6.904	99	1187920	37.312	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.063	67	769586	36.961	ng/ul	0.00
11) 2-Chlorophenol-d4	7.257	132	1026050	38.845	ng/ul	0.00
15) 4-Methylphenol-d8	8.440	113	957515	38.045	ng/ul	0.00
21) Nitrobenzene-d5	8.881	128	466504	45.204	ng/ul	0.00
24) 2-Nitrophenol-d4	9.598	143	460711	53.960	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	796572	37.759	ng/ul	0.00
31) 4-Chloroaniline-d4	10.663	131	1303635	37.856	ng/ul	0.00
46) Dimethylphthalate-d6	13.798	166	2290521	38.686	ng/ul	0.00
49) Acenaphthylene-d8	14.069	160	2800552	37.061	ng/ul	0.00
54) 4-Nitrophenol-d4	14.598	143	323132m	34.941	ng/ul	0.02
60) Fluorene-d10	15.374	176	1894144	37.279	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	157671	30.993	ng/ul	0.00
73) Anthracene-d10	17.239	188	2709705	39.664	ng/ul	0.00
81) Pyrene-d10	19.486	212	2882125	37.258	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.468	264	3102231	39.791	ng/ul	0.00

Target Compounds Qvalue

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

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