

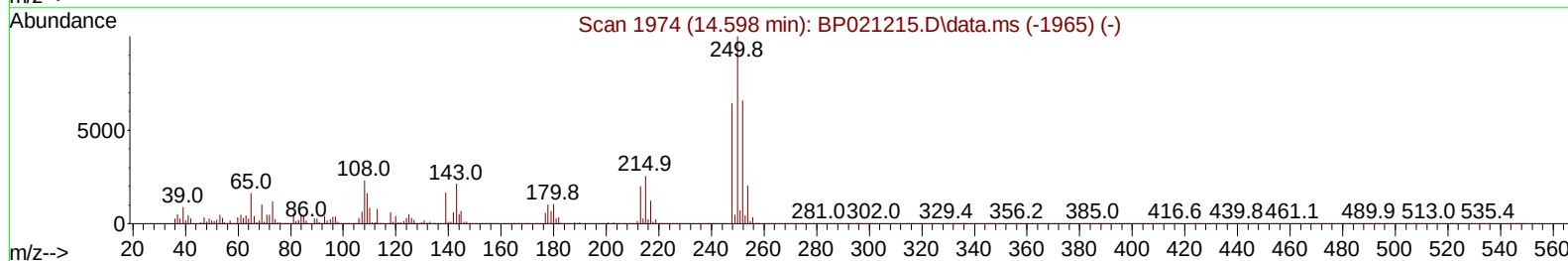
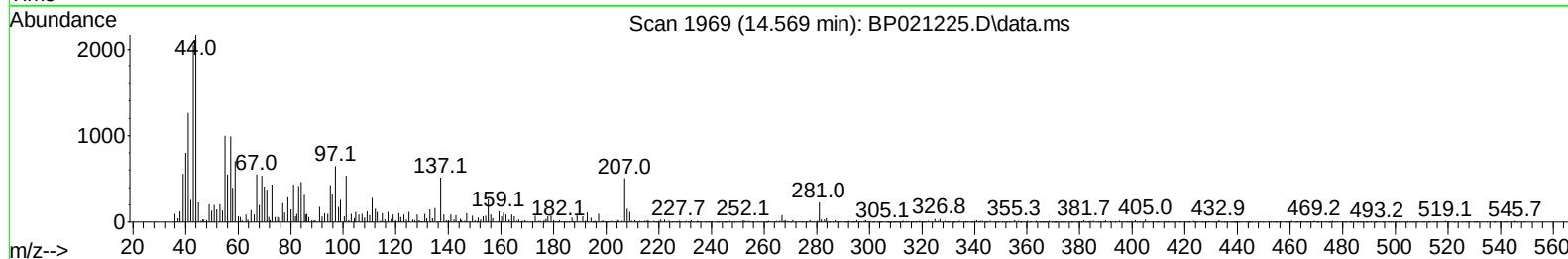
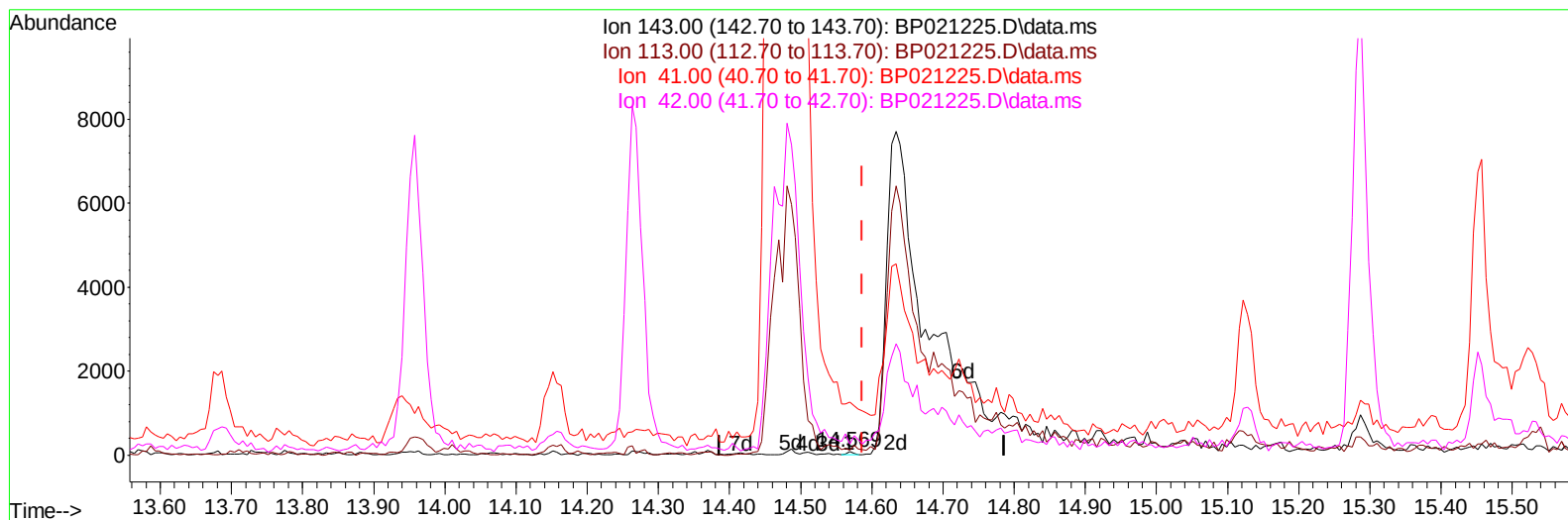
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021225.D
Acq On : 30 Jul 2024 02:34
Operator : MA/JU
Sample : P3300-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G92

Manual IntegrationsAPPROVED

Quant Time: Jul 30 03:43:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/30/2024
Supervised By :mohammad ahmed 09/04/2024



TIC: BP021225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.017) 0.01 ng/u1

response 36

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	143.75#
41.00	37.90	1573.75#
42.00	27.10	323.75#

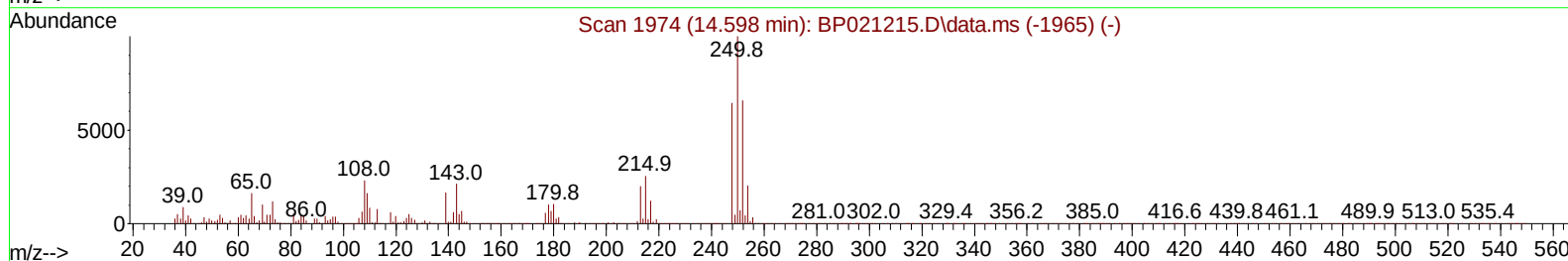
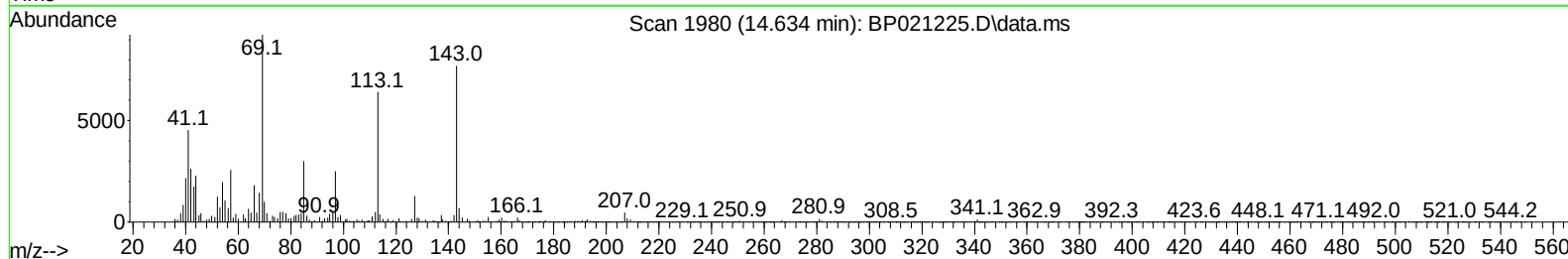
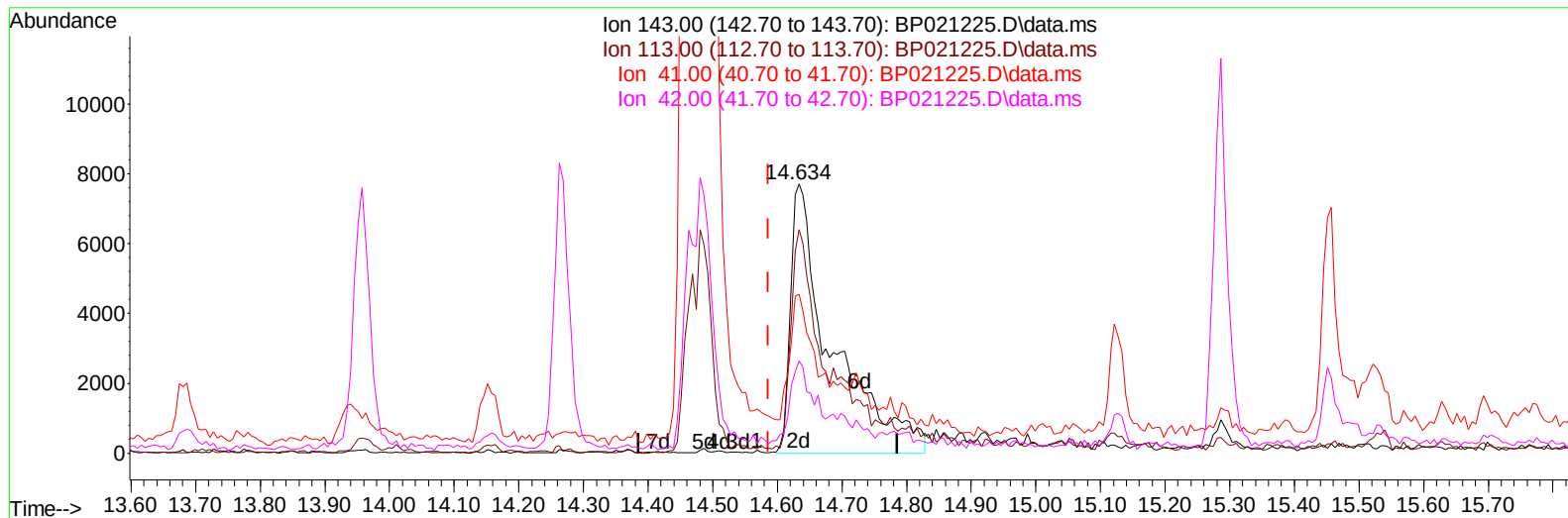
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TIC: BP021225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.047) 7.42 ng/ul m

response 34028

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	83.10
41.00	37.90	59.09#
42.00	27.10	34.30#

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Instrument :
 BNA_P
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 E1G92

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/30/2024
 Supervised By :mohammad ahmed 09/04/2024

Quant Time: Jul 30 03:46:07 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.634	152		186984	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136		709843	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164		443131	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188		946157	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240		1022253	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264		1224145	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.170	96		18015	4.384	ng/uL	0.00
4) Pyridine-d5	3.605	84		78951	6.637	ng/ul	0.01
7) Phenol-d5	6.864	99		109882	7.208	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67		229909	24.886	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132		289487	23.048	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113		203123	16.043	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128		139216	25.250	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143		158059	25.047	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.052	165		286706	25.126	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131		366281	24.230	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166		951331	29.532	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160		1012981	28.005	ng/ul	0.00
54) 4-Nitrophenol-d4	14.634	143		34028m	7.424	ng/ul	0.05
60) Fluorene-d10	15.287	176		725368	27.447	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200		138503	24.194	ng/ul	0.00
73) Anthracene-d10	17.192	188		1352879	32.193	ng/ul	0.00
81) Pyrene-d10	19.575	212		1687464	26.700	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264		2020576	33.467	ng/ul	0.00
Target Compounds							
86) Bis(2-ethylhexyl)phtha...	21.422	149		126583	2.718	ng/ul	Qvalue 99

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

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