

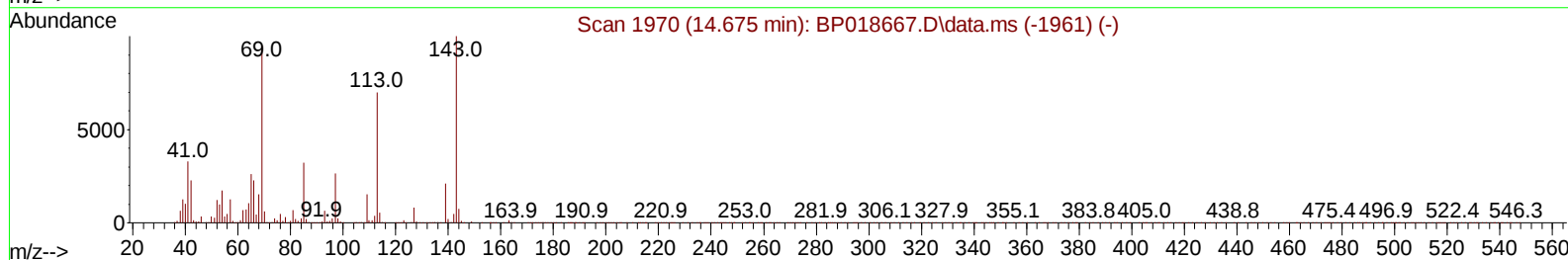
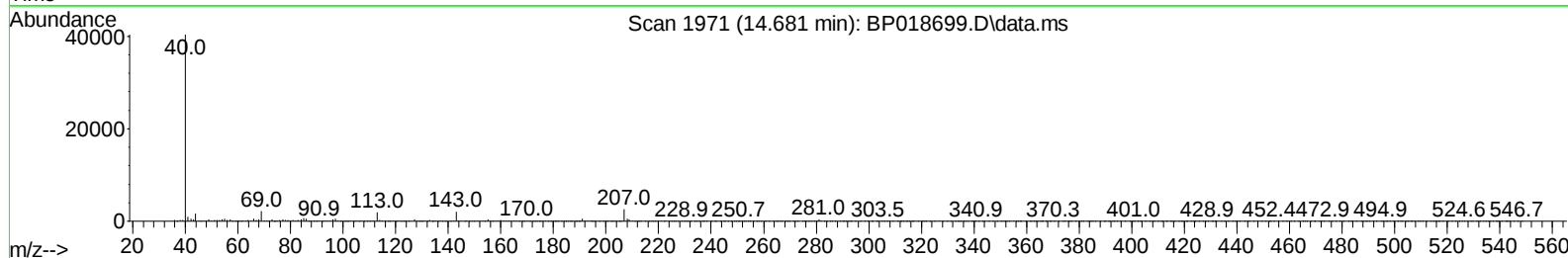
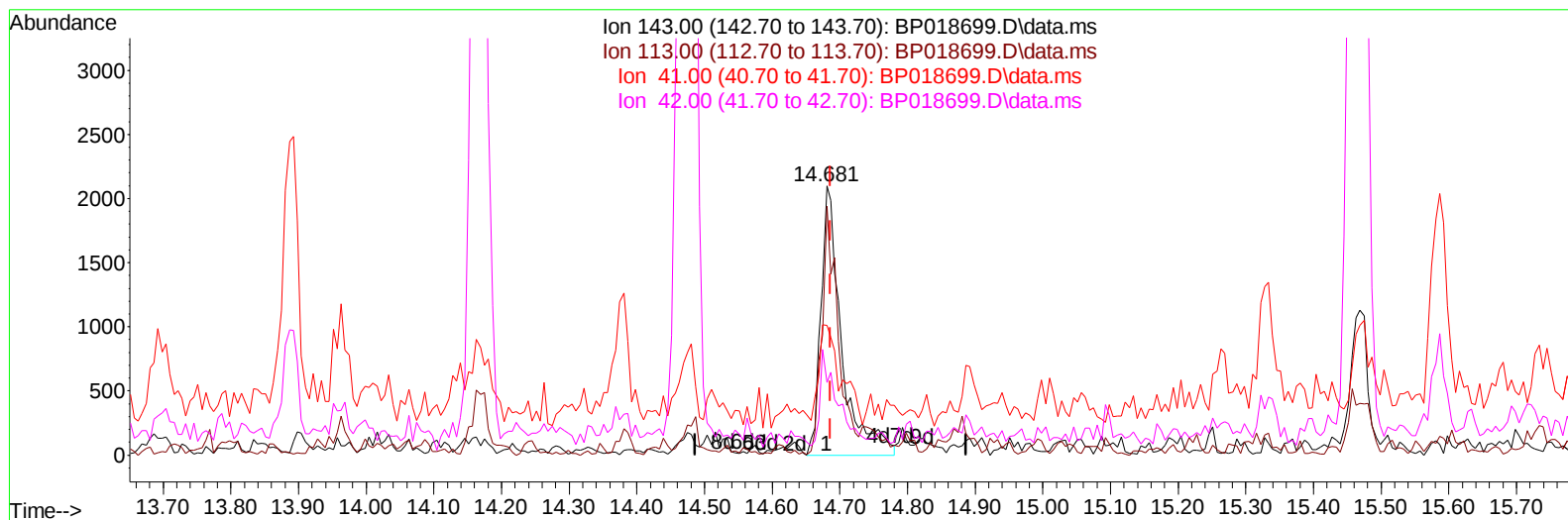
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120723\
 Data File : BP018699.D
 Acq On : 08 Dec 2023 06:27
 Operator : MA/JU
 Sample : 05497-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB0

Manual IntegrationsAPPROVED

Quant Time: Dec 08 07:04:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 04 21:39:40 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/08/2023
 Supervised By :mohammad ahmed 12/08/2023



TIC: BP018699.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.006) 0.41 ng/ul m

response 4470

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	92.47#
41.00	34.60	48.16#
42.00	23.00	26.94

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Quant Time: Dec 08 07:05:46 2023
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.846	152		380205	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.640	136		1484121	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.475	164		740387	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.222	188		2208247	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240		2383528	20.000	ng/ul	0.00
88) Perylene-d12	23.680	264		2762971	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		32463	2.910	ng/uL	0.00
4) Pyridine-d5	3.681	84		379562	12.701	ng/ul	0.00
7) Phenol-d5	7.016	99		155517	4.076	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67		317734	14.050	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.375	132		214850	7.216	ng/ul	-0.01
15) 4-Methylphenol-d8	8.557	113		58455	1.893	ng/ul	-0.01
21) Nitrobenzene-d5	9.004	128		198740	15.046	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.728	143		198935	14.526	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.263	165		230028	8.214	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.781	131		257845	7.046	ng/ul	-0.01
46) Dimethylphthalate-d6	13.892	166		1199065	18.146	ng/ul	-0.01
49) Acenaphthylene-d8	14.169	160		1123548	15.441	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143		4470m	0.412	ng/ul	0.00
60) Fluorene-d10	15.469	176		1096941	19.776	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200		37663	2.844	ng/ul	0.00
73) Anthracene-d10	17.322	188		1730377	14.905	ng/ul	0.00
81) Pyrene-d10	19.557	212		2252394	12.140	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.527	264		1914631	11.885	ng/ul	0.00

Target Compounds Qvalue

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

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