

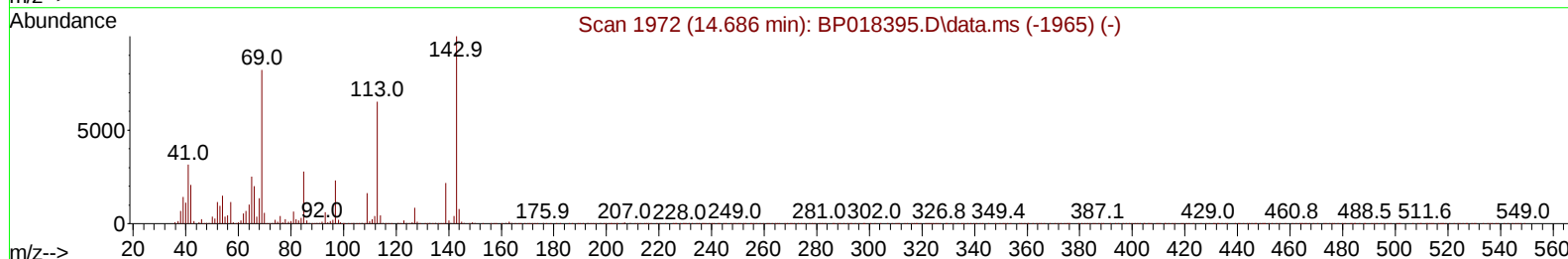
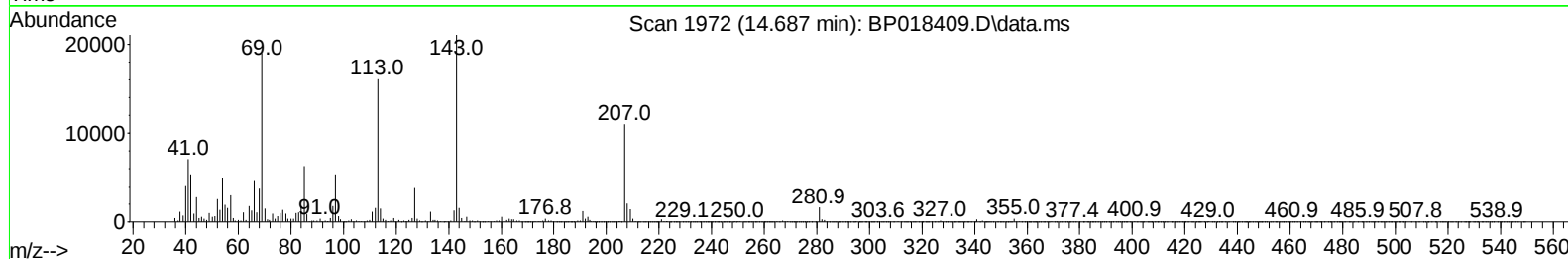
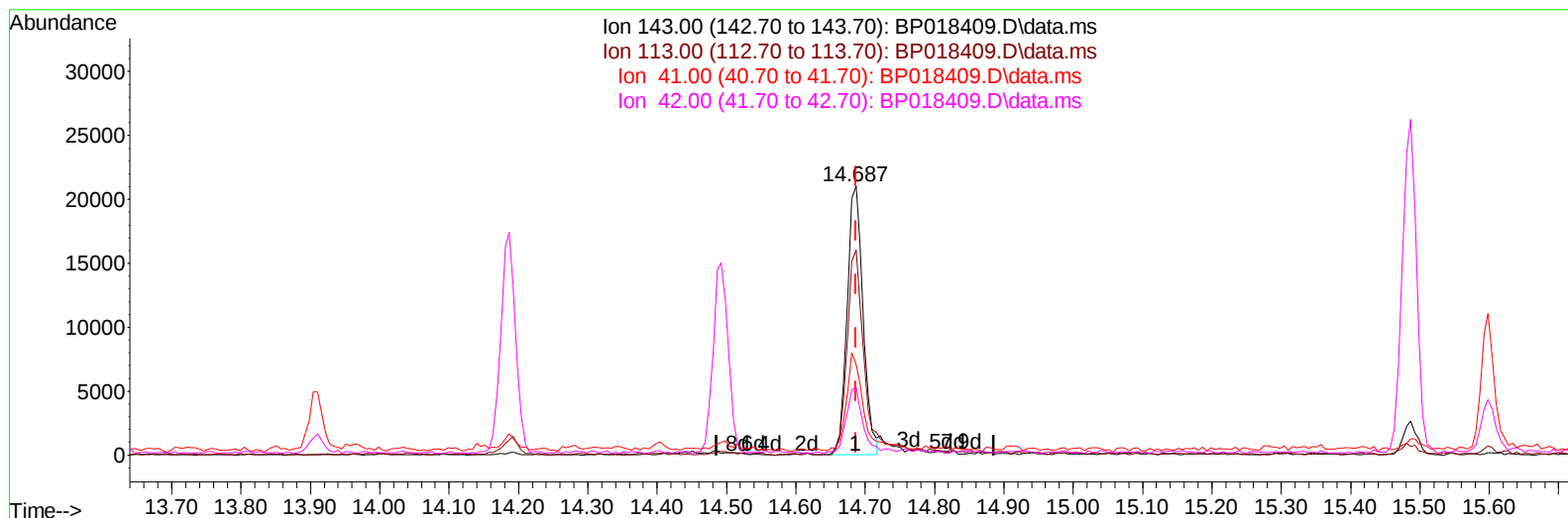
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018409.D
 Acq On : 27 Nov 2023 20:03
 Operator : MA/JU
 Sample : 05332-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG64

Manual IntegrationsAPPROVED

Quant Time: Nov 27 21:58:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 27 21:50:01 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/28/2023
 Supervised By :mohammad ahmed 12/01/2023



TIC: BP018409.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.000) 2.67 ng/u1

response 33324

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.70	76.39
41.00	36.10	33.54
42.00	25.00	25.47

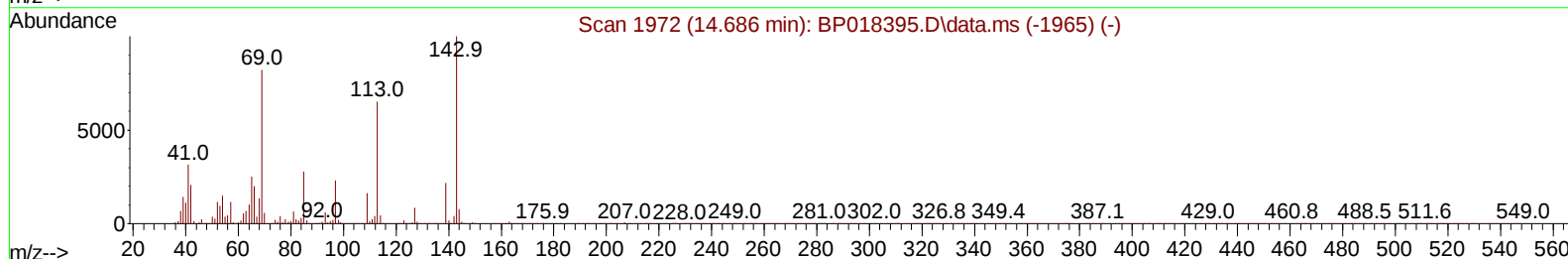
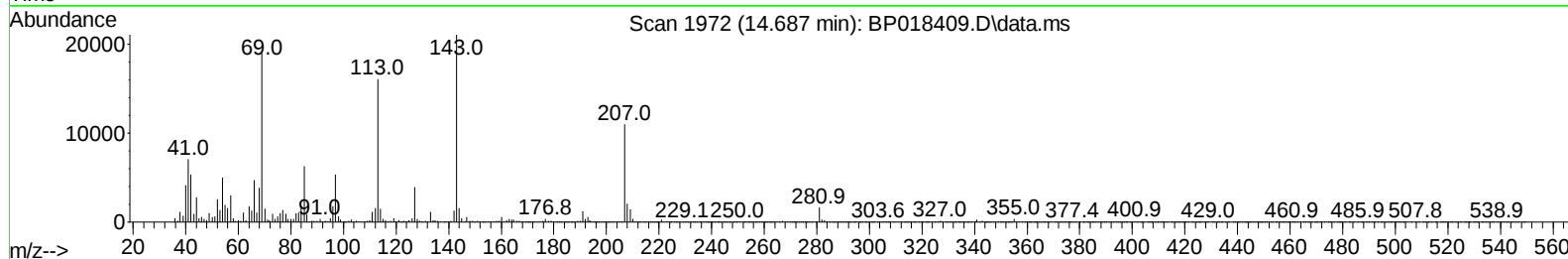
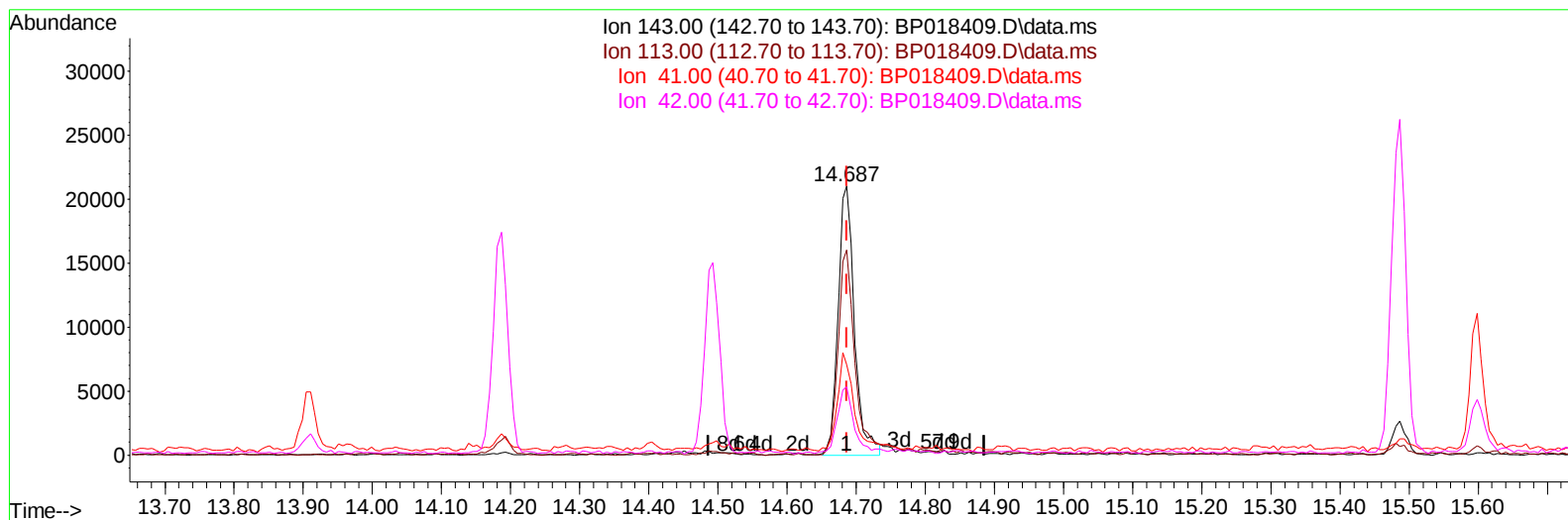
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TIC: BP018409.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.000) 2.77 ng/ul m

response 34537

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.70	76.39
41.00	36.10	33.54
42.00	25.00	25.47

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/28/2023
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Quant Time: Nov 27 22:28:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 27 21:50:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	444161	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1778035	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	1014629	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1778183	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	1466687	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	1726272	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	40632	3.426	ng/uL	0.00
4) Pyridine-d5	3.687	84	248773	7.634	ng/ul	0.00
7) Phenol-d5	7.022	99	215683	5.313	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	656747	27.057	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	696830	20.446	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	420887	12.848	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	442843	30.664	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	396762	29.748	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	779590	24.227	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	936294	21.574	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2573610	28.506	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	2909978	28.288	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	34537m	2.771	ng/ul	0.00
60) Fluorene-d10	15.487	176	2129106	28.195	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	201906	26.673	ng/ul	0.00
73) Anthracene-d10	17.345	188	3259161	34.291	ng/ul	0.00
81) Pyrene-d10	19.592	212	3413748	29.245	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	3707967	37.049	ng/ul	0.03

Target Compounds Qvalue

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

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