

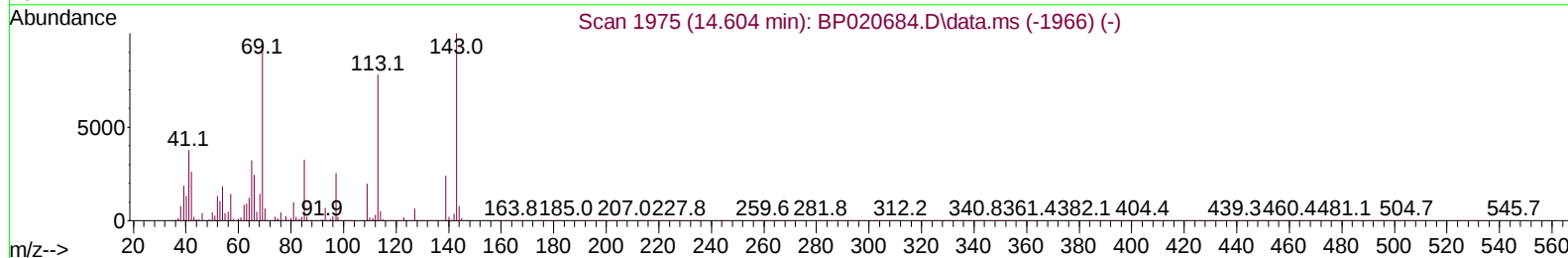
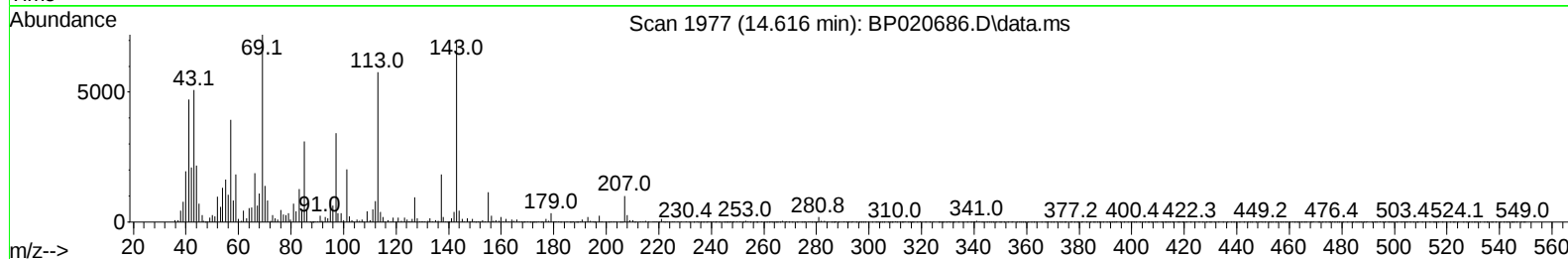
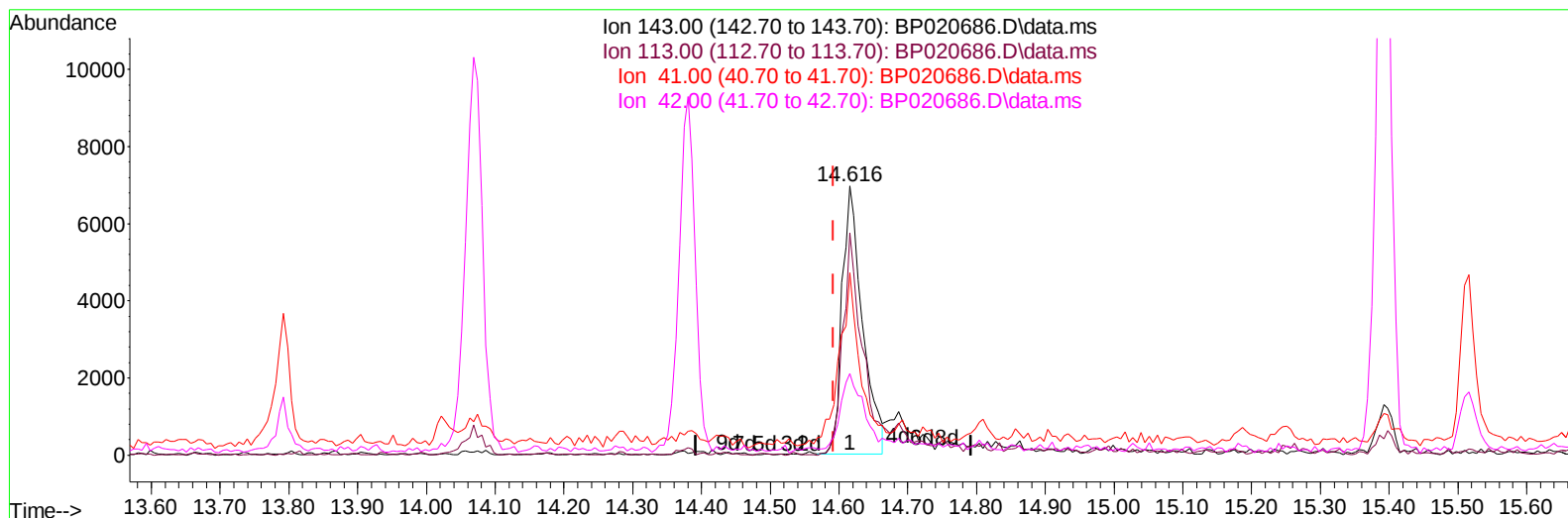
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020686.D
Acq On : 28 Jun 2024 09:47
Operator : MA/JU
Sample : P2898-01
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G63

Manual IntegrationsAPPROVED

Quant Time: Jun 28 10:51:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 25 18:30:06 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/28/2024
Supervised By :mohammad ahmed 06/29/2024



TIC: BP020686.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.023) 2.66 ng/u1

response 14282

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.40	82.59
41.00	40.20	67.82#
42.00	28.00	30.30

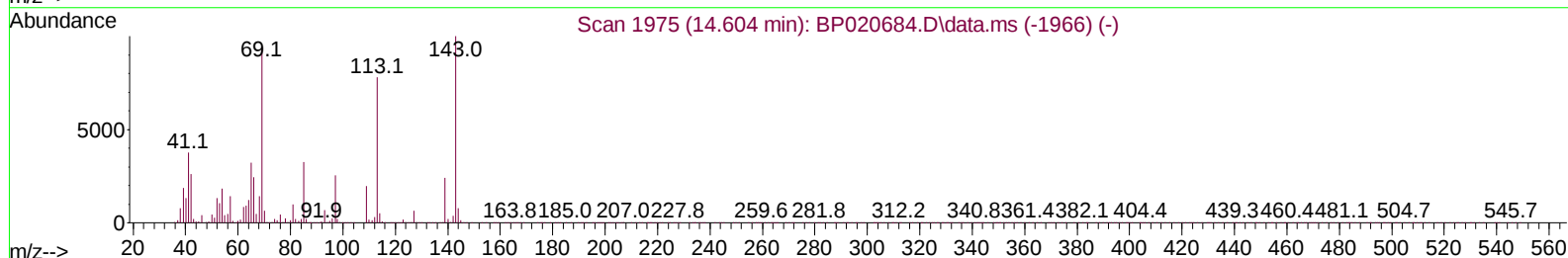
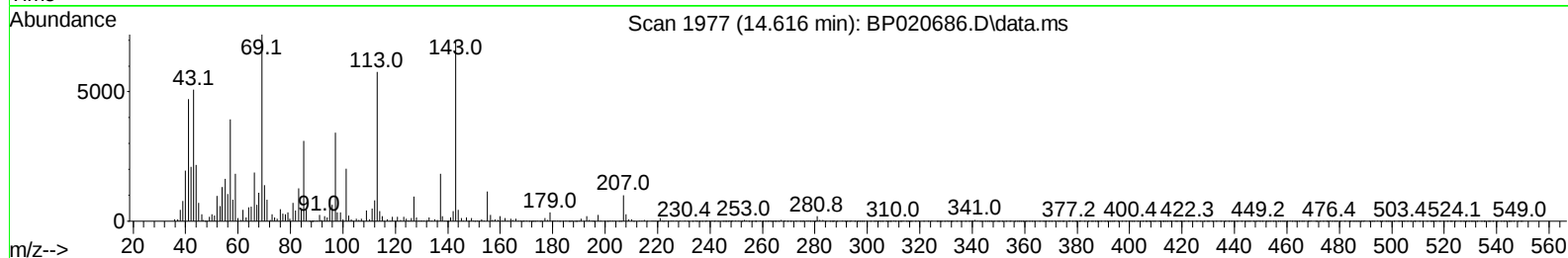
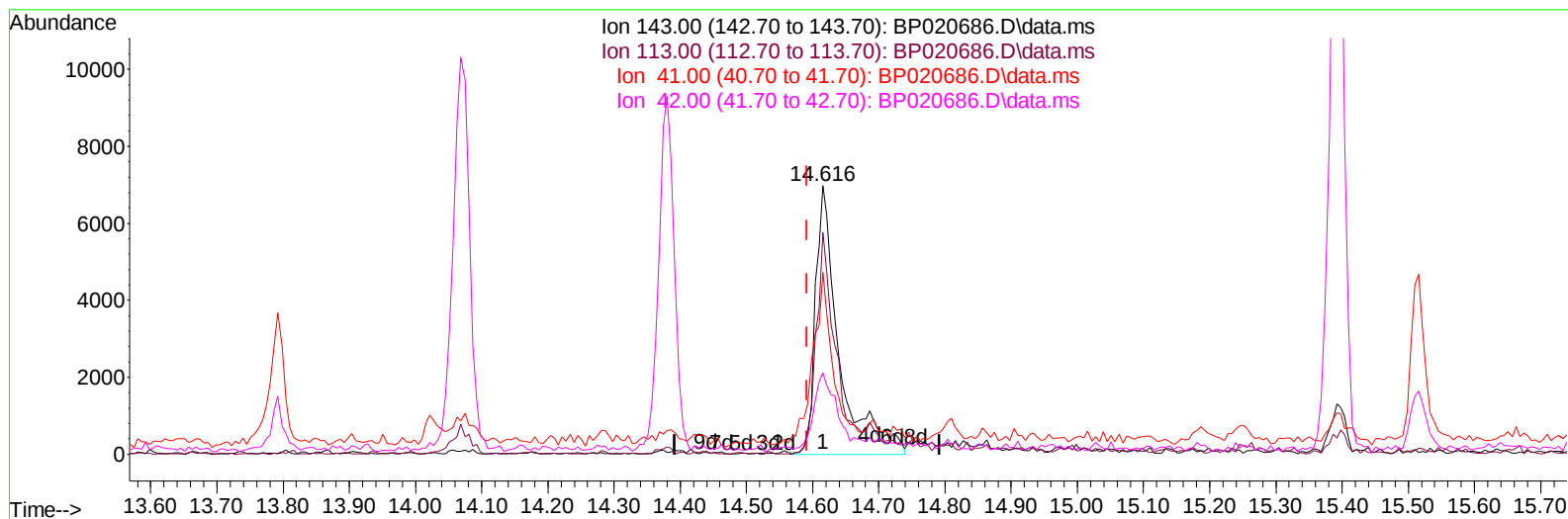
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TIC: BP020686.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.023) 3.24 ng/ul m

response 17403

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.40	82.59
41.00	40.20	67.82#
42.00	28.00	30.30

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 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1G63

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/28/2024
 Supervised By :mohammad ahmed 06/29/2024

Quant Time: Jun 28 10:53:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 25 18:30:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.745	152	210909	20.000	ng/ul	0.00
20) Naphthalene-d8	10.516	136	835583	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.380	164	505579	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.192	188	903292	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240	716969	20.000	ng/ul	0.01
88) Perylene-d12	25.027	264	911892	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	20480	4.170	ng/uL	0.00
4) Pyridine-d5	3.693	84	133599	9.333	ng/ul	0.00
7) Phenol-d5	6.922	99	84395	4.879	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.087	67	322326	29.392	ng/ul	0.00
11) 2-Chlorophenol-d4	7.287	132	225209	15.930	ng/ul	0.00
15) 4-Methylphenol-d8	8.440	113	187639	13.356	ng/ul	0.00
21) Nitrobenzene-d5	8.893	128	220339	35.717	ng/ul	0.00
24) 2-Nitrophenol-d4	9.598	143	146190	23.680	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.145	165	249908	19.652	ng/ul	0.00
31) 4-Chloroaniline-d4	10.657	131	487521	27.356	ng/ul	0.00
46) Dimethylphthalate-d6	13.792	166	1383771	39.365	ng/ul	0.01
49) Acenaphthylene-d8	14.069	160	1497462	35.700	ng/ul	0.00
54) 4-Nitrophenol-d4	14.616	143	17403m	3.240	ng/ul	0.02
60) Fluorene-d10	15.392	176	1102661	37.277	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	77426	17.009	ng/ul	0.01
73) Anthracene-d10	17.292	188	1614171	39.817	ng/ul	0.01
81) Pyrene-d10	19.674	212	1650570	38.321	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.803	264	1788388	40.317	ng/ul	0.04

Target Compounds Qvalue

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

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