

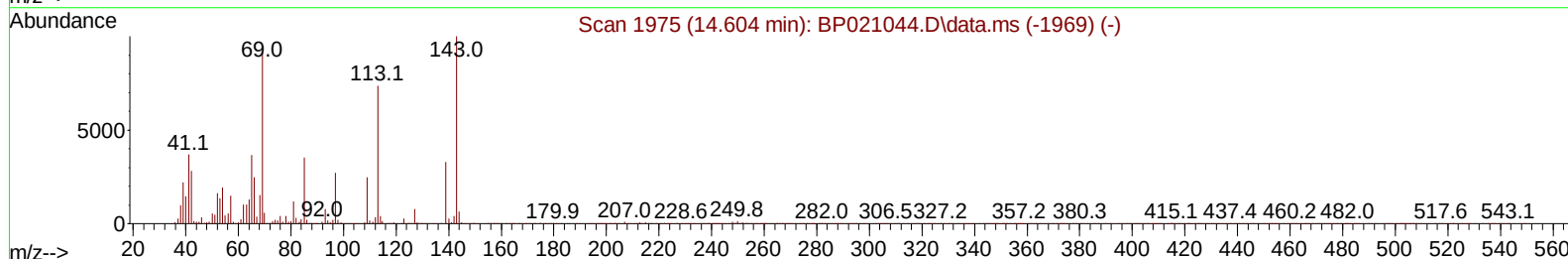
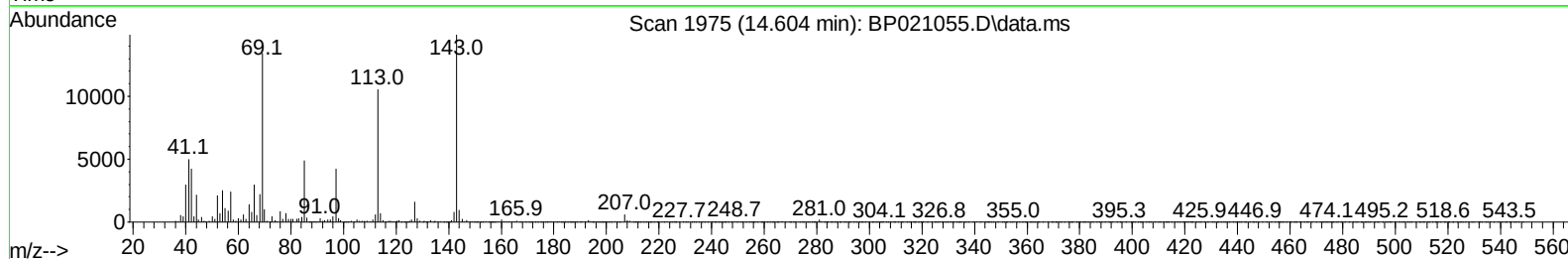
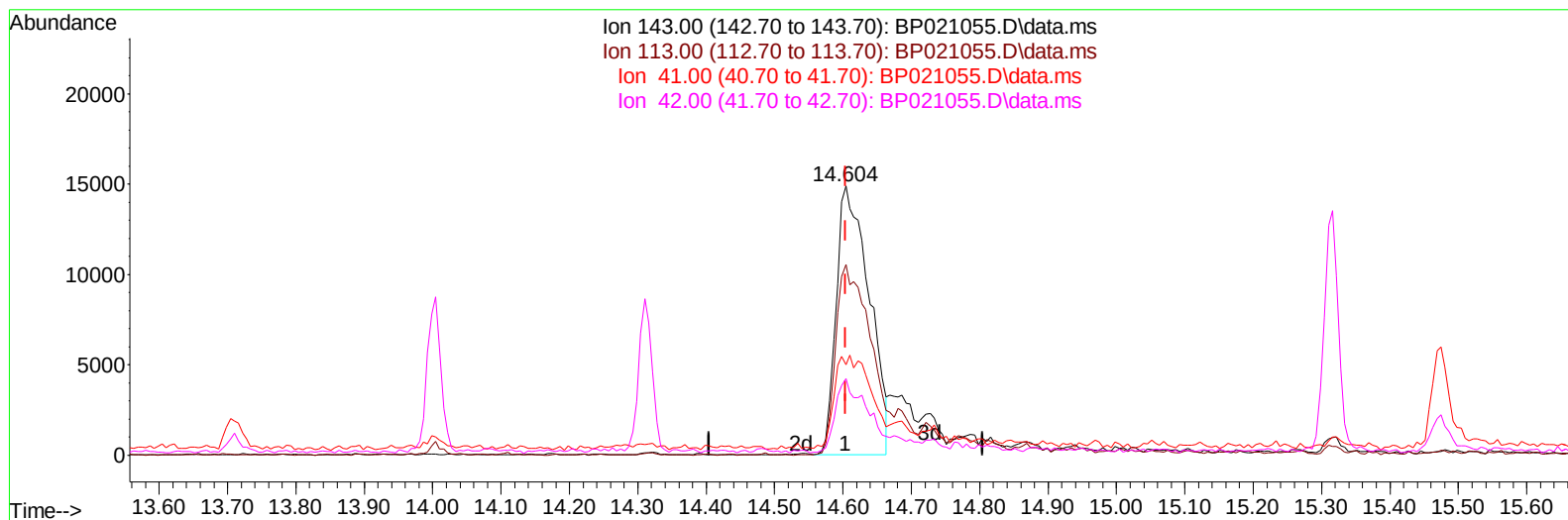
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021055.D
Acq On : 18 Jul 2024 21:13
Operator : MA/JU
Sample : P3260-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E08H0

Manual IntegrationsAPPROVED

Quant Time: Jul 18 22:56:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 18 14:23:21 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/19/2024
Supervised By :mohammad ahmed 07/20/2024



TIC: BP021055.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.604min (-0.000) 11.83 ng/u1

response 49583

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	70.97
41.00	37.90	33.72
42.00	27.10	28.40

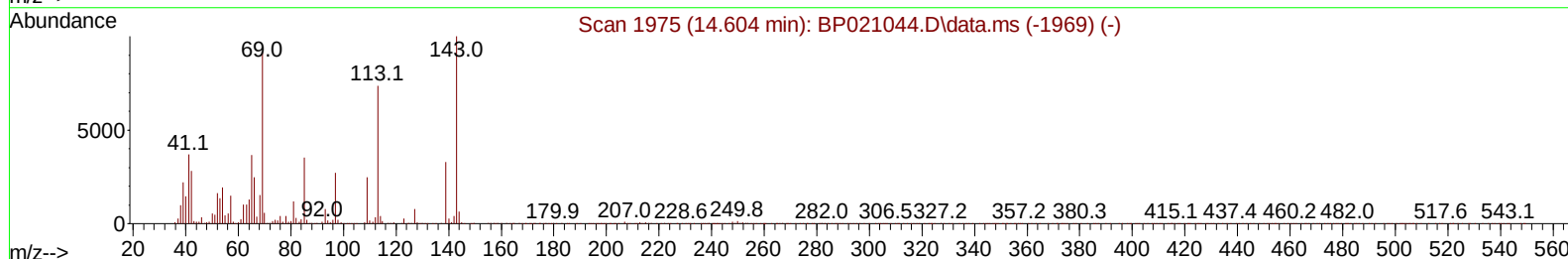
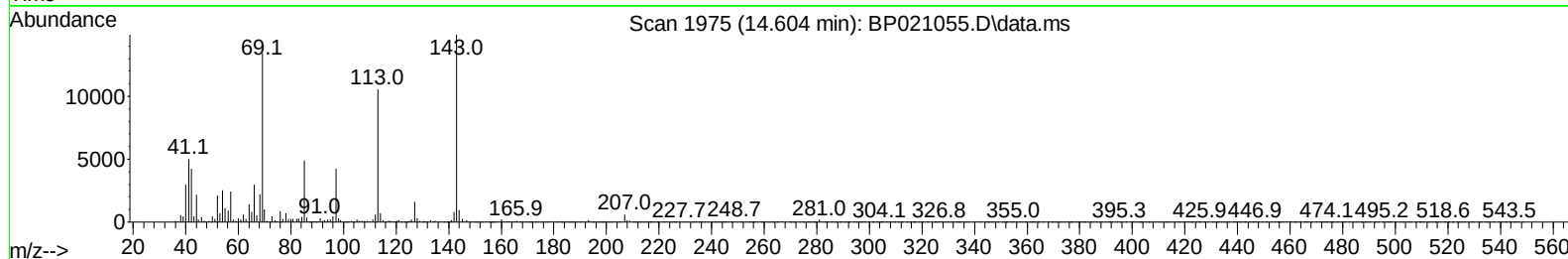
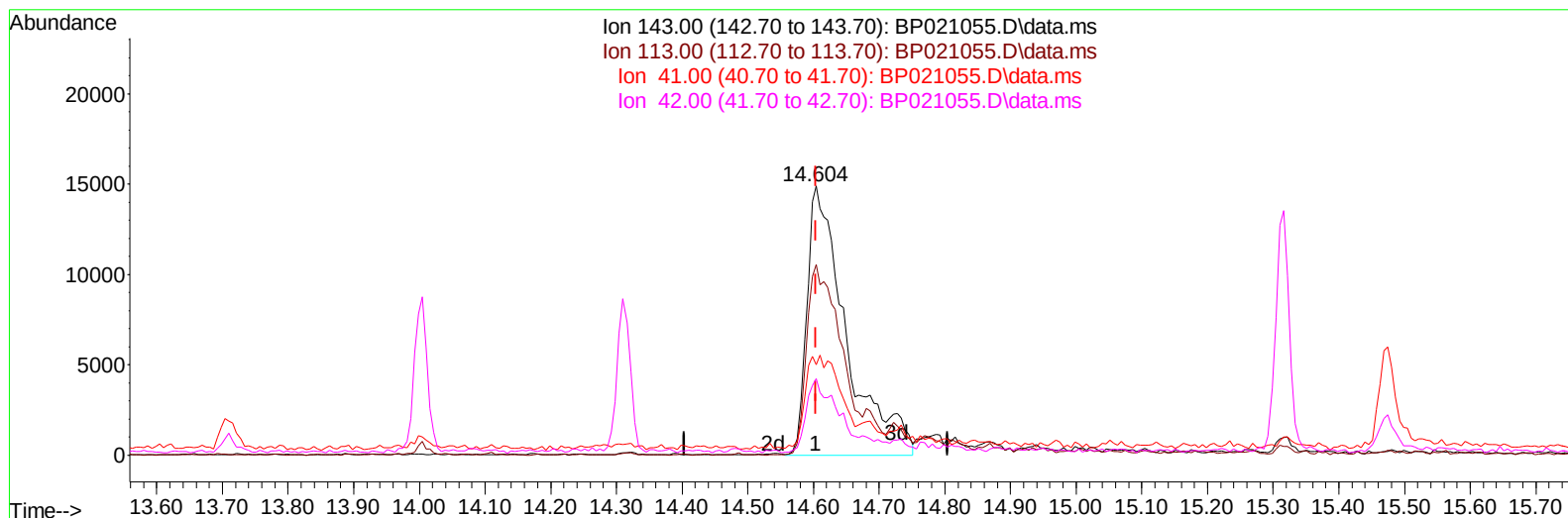
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TIC: BP021055.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.604min (-0.000) 14.74 ng/ul m

response 61747

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	70.97
41.00	37.90	33.72
42.00	27.10	28.40

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 Misc :
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Instrument :
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 ClientSampleId :
 E08H0

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/19/2024
 Supervised By :mohammad ahmed 07/20/2024

Quant Time: Jul 18 23:19:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 18 14:23:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	162617	20.000	ng/ul	0.00
20) Naphthalene-d8	10.451	136	637552	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	405079	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.122	188	887374	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	862053	20.000	ng/ul	0.00
88) Perylene-d12	24.880	264	1022323	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	21919	6.134	ng/uL	0.00
4) Pyridine-d5	3.652	84	126076	12.186	ng/ul	0.00
7) Phenol-d5	6.899	99	186597	14.075	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	243255	30.276	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	331226	30.323	ng/ul	0.00
15) 4-Methylphenol-d8	8.404	113	265654	24.126	ng/ul	-0.01
21) Nitrobenzene-d5	8.834	128	162026	32.720	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.546	143	184110	32.484	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	317376	30.968	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.598	131	372670	27.448	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	968280	32.882	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1123465	33.977	ng/ul	0.00
54) 4-Nitrophenol-d4	14.604	143	61747m	14.738	ng/ul	0.00
60) Fluorene-d10	15.316	176	837398	34.663	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	135512	25.239	ng/ul	0.00
73) Anthracene-d10	17.222	188	1395856	35.416	ng/ul	0.00
81) Pyrene-d10	19.604	212	1725013	32.367	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.657	264	1842418	36.541	ng/ul	-0.03
Target Compounds						
					Qvalue	
6) Benzaldehyde	6.863	77	15283	2.287	ng/ul	96

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

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