

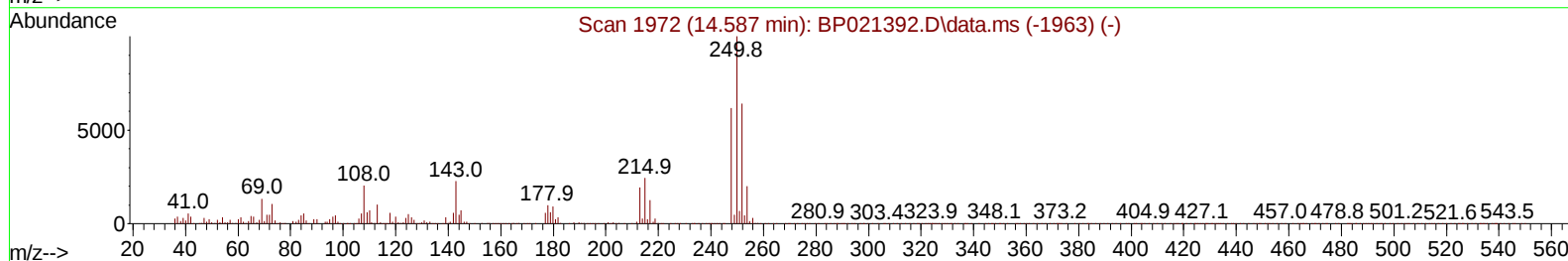
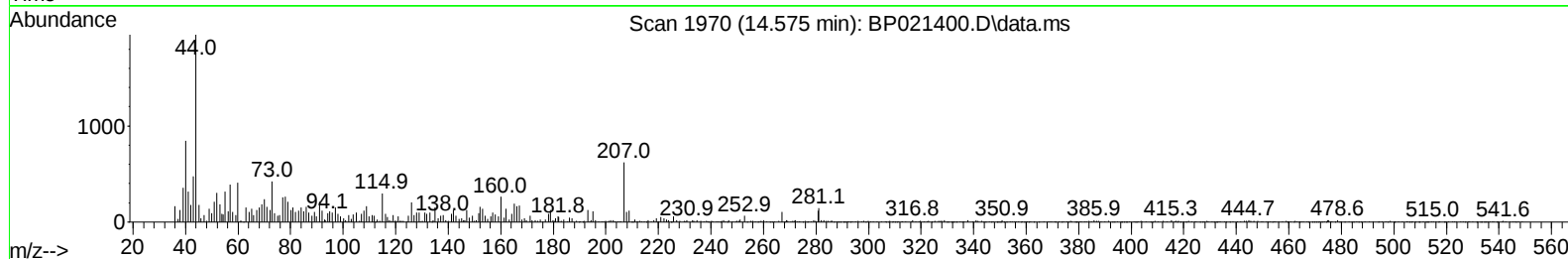
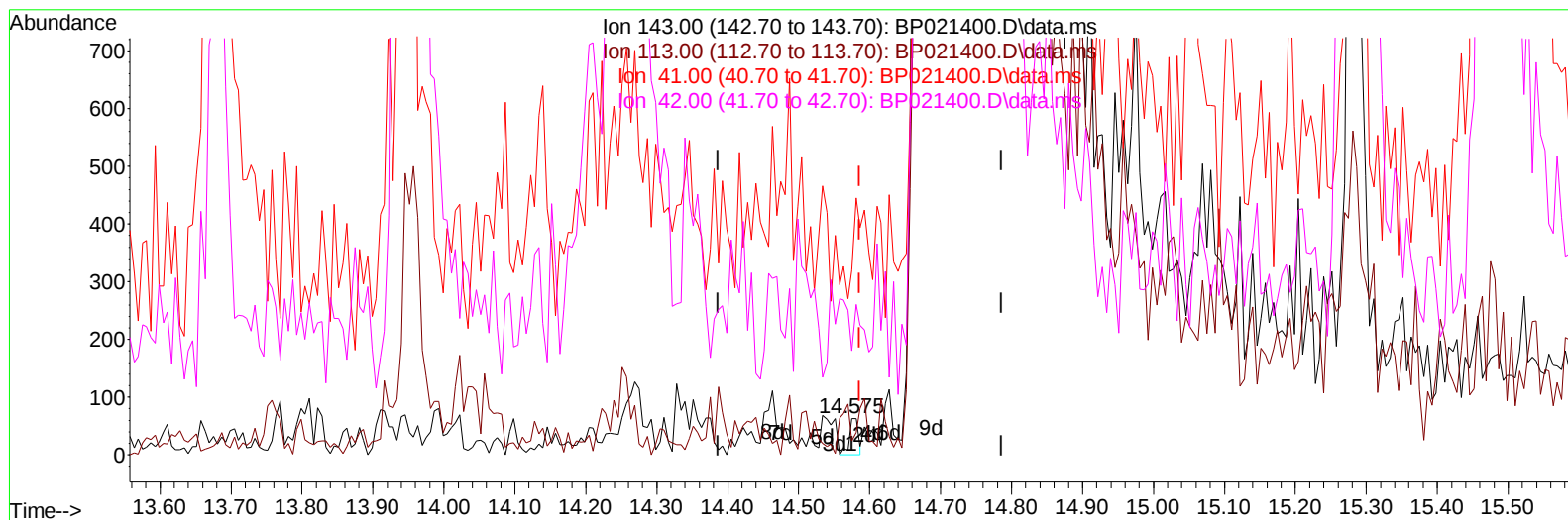
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
Data File : BP021400.D
Acq On : 06 Aug 2024 21:41
Operator : CG/JU
Sample : P3378-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY3

Manual IntegrationsAPPROVED

Quant Time: Aug 06 22:52:37 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 08/07/2024
Supervised By :mohammad ahmed 08/09/2024



TIC: BP021400.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (-0.011) 0.01 ng/u1

response 56

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	38.10#
41.00	37.90	507.94#
42.00	27.10	285.71#

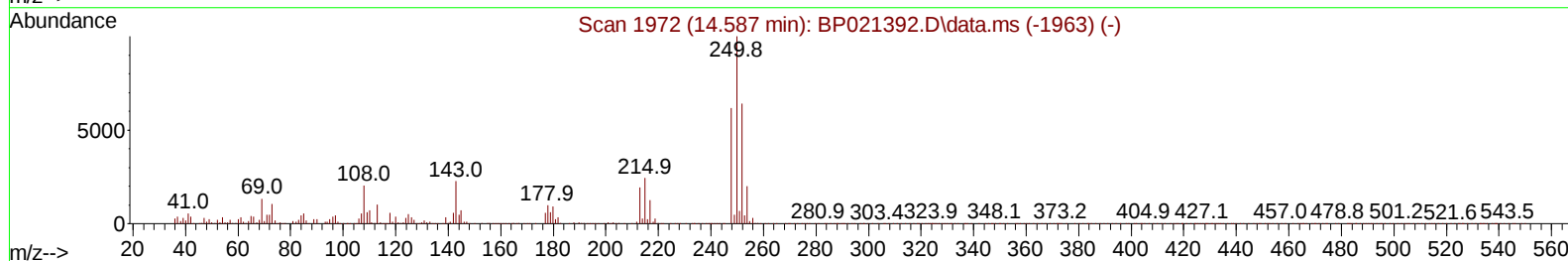
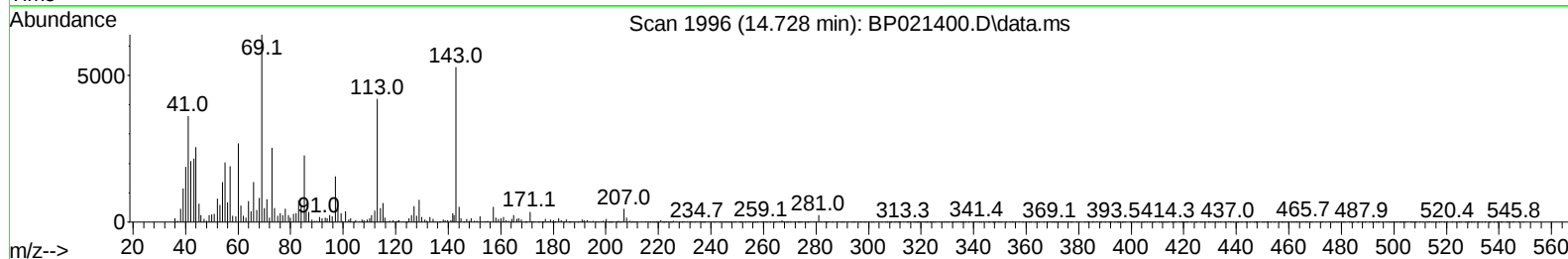
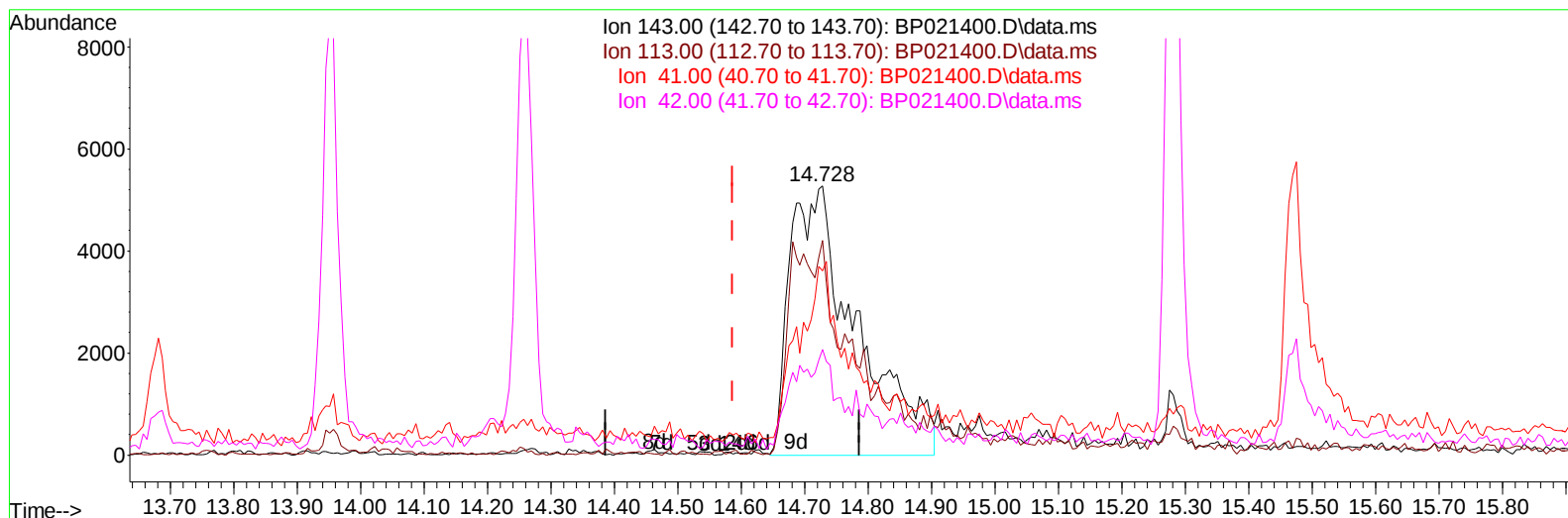
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TIC: BP021400.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.142) 7.01 ng/ul m

response 38653

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.81
41.00	37.90	68.52#
42.00	27.10	39.35#

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 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/07/2024
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Quant Time: Aug 06 23:27:17 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.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	186059	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.387	136	764480	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	533474	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.087	188	1202357	20.000	ng/ul	0.00
79) Chrysene-d12	21.528	240	1286210	20.000	ng/ul	0.00
88) Perylene-d12	24.810	264	1530565	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	17624	4.310	ng/uL	-0.02
4) Pyridine-d5	3.593	84	93050	7.861	ng/ul	0.00
7) Phenol-d5	6.875	99	85155	5.614	ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	6.969	67	239162	26.016	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.175	132	284696	22.780	ng/ul	0.00
15) 4-Methylphenol-d8	8.364	113	198265	15.737	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	161836	27.255	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.493	143	186564	27.451	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.063	165	305366	24.849	ng/ul	0.01
31) 4-Chloroaniline-d4	10.563	131	389915	23.950	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	1167777	30.112	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1257333	28.874	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.728	143	38653m	7.005	ng/ul	0.14
60) Fluorene-d10	15.281	176	993077	31.214	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	161513	22.201	ng/ul	0.01
73) Anthracene-d10	17.187	188	1636876	30.652	ng/ul	0.00
81) Pyrene-d10	19.563	212	2149944	27.037	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.569	264	2472550	32.755	ng/ul	-0.03

Target Compounds Qvalue

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

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