Quantitation Report (Qedit)

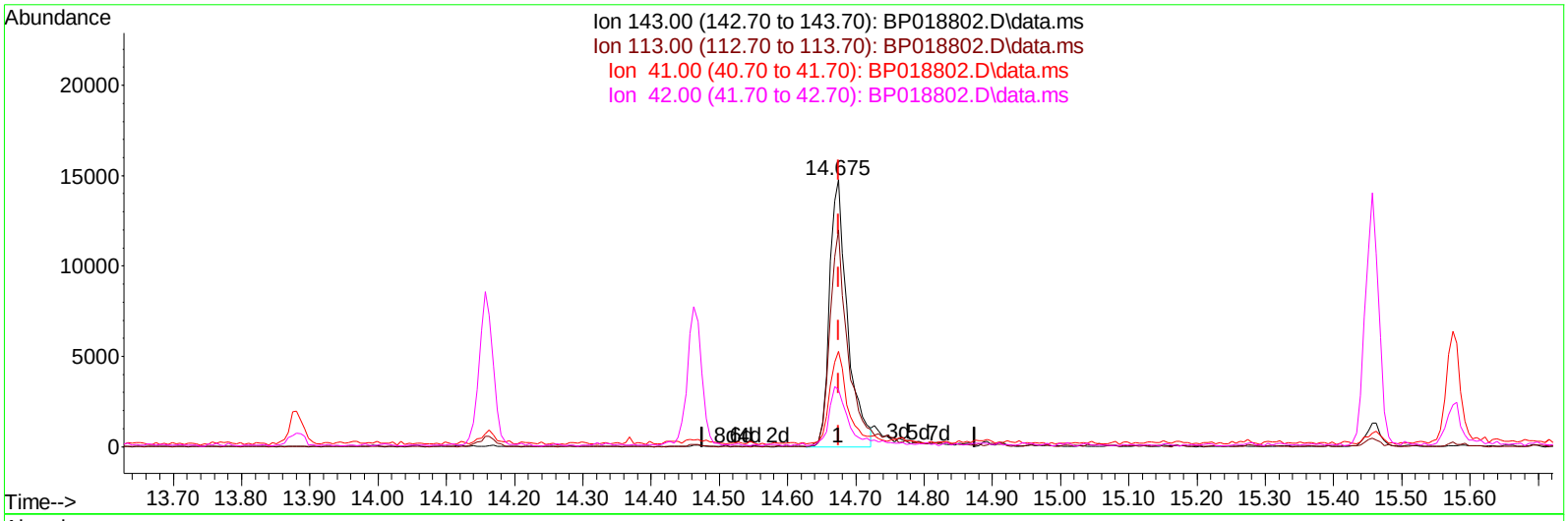
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
Data File : BP018802.D
Acq On : 13 Dec 2023 11:31
Operator : MA/JU
Sample : 05673-06
Misc :
ALS Vial : 82 Sample Multiplier: 1

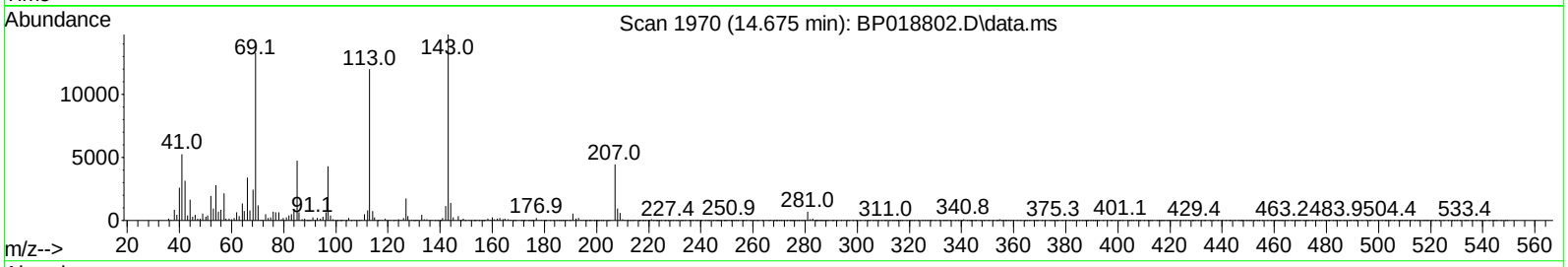
Quant Time: Dec 13 21:36:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 11 22:02:13 2023
Response via : Initial Calibration

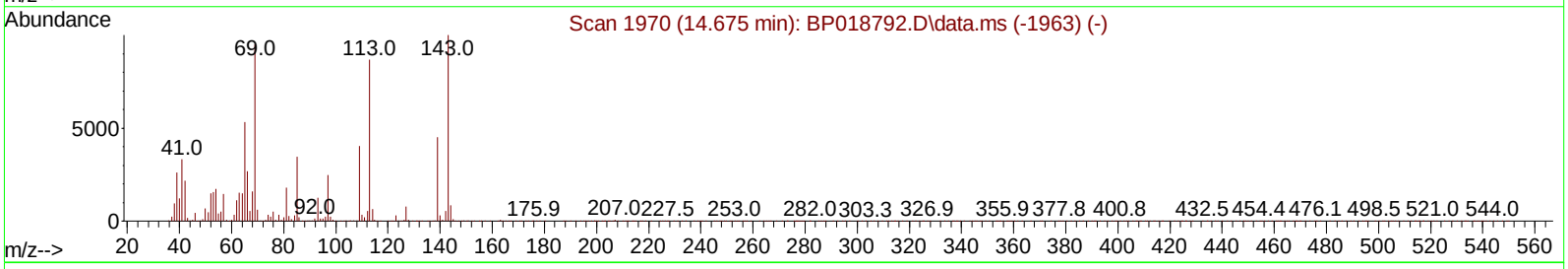
Instrument :
BNA_P
ClientSampleId :
BHEH9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/14/2023
Supervised By :mohammad ahmed 12/14/2023







TIC: BP018802.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (-0.000) 4.42 ng/ul

response 27090

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	81.20
41.00	34.60	35.77
42.00	23.00	21.34

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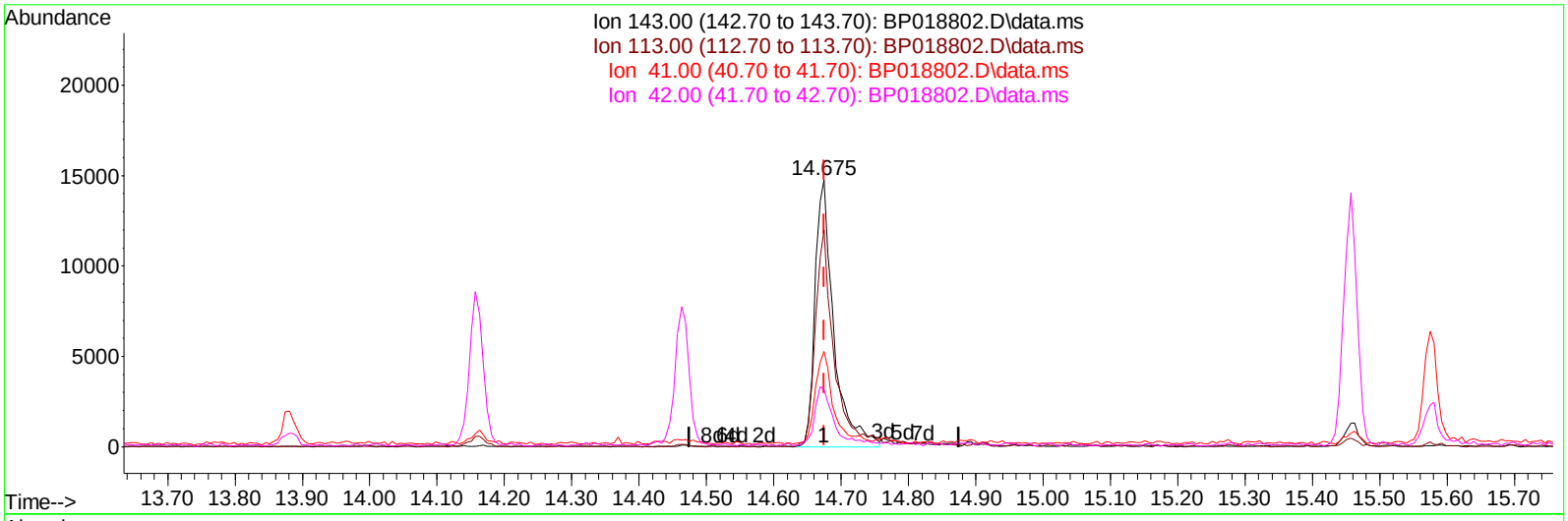

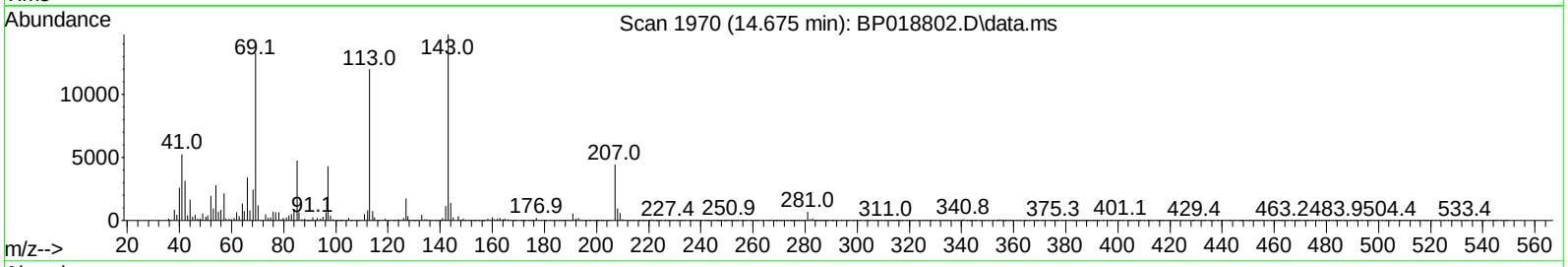

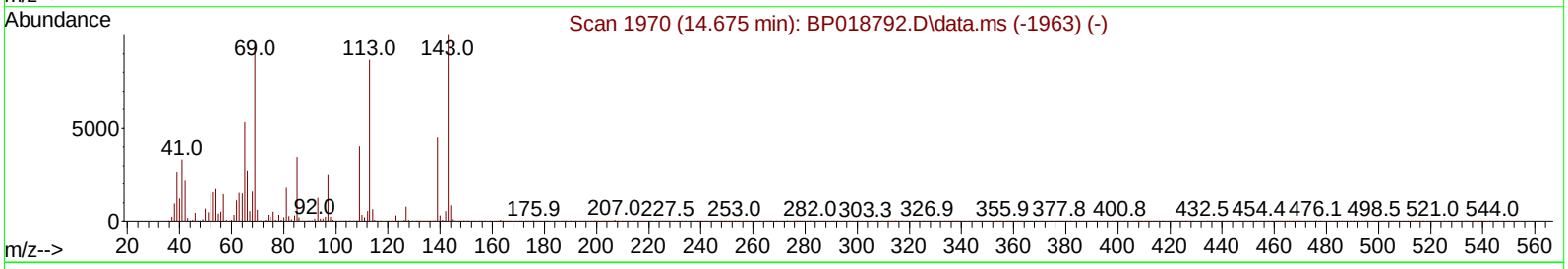

TIC: BP018802.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (-0.000) 4.64 ng/ul m

response 28467

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	81.20
41.00	34.60	35.77
42.00	23.00	21.34

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
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Acq On : 13 Dec 2023 11:31
Operator : MA/JU
Sample : 05673-06
Misc :
ALS Vial : 82 Sample Multiplier: 1

Quant Time: Dec 13 21:47:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 11 22:02:13 2023
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	172221	20.000	ng/ul	0.00
20) Naphthalene-d8	10.628	136	649190	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	418579	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1001419	20.000	ng/ul	-0.01
79) Chrysene-d12	21.298	240	1093836	20.000	ng/ul	-0.01
88) Perylene-d12	23.662	264	1273802	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	20755	4.107	ng/uL	0.00
4) Pyridine-d5	3.687	84	136540	10.086	ng/ul	0.01
7) Phenol-d5	7.004	99	93305	5.399	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	265375	25.907	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	271893	20.161	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	174433	12.471	ng/ul	0.00
21) Nitrobenzene-d5	8.998	128	168309	29.129	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	170116	28.398	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	318683	26.017	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	414803	25.912	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1095490	29.324	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1173132	28.518	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	28467m	4.642	ng/ul	0.00
60) Fluorene-d10	15.457	176	949170	30.268	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	145905	24.299	ng/ul	-0.01
73) Anthracene-d10	17.310	188	1657439	31.482	ng/ul	-0.01
81) Pyrene-d10	19.551	212	2116467	24.857	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.509	264	2324569	31.298	ng/ul	-0.02

Target Compounds Qvalue

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

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