Quantitation Report (Qedit)

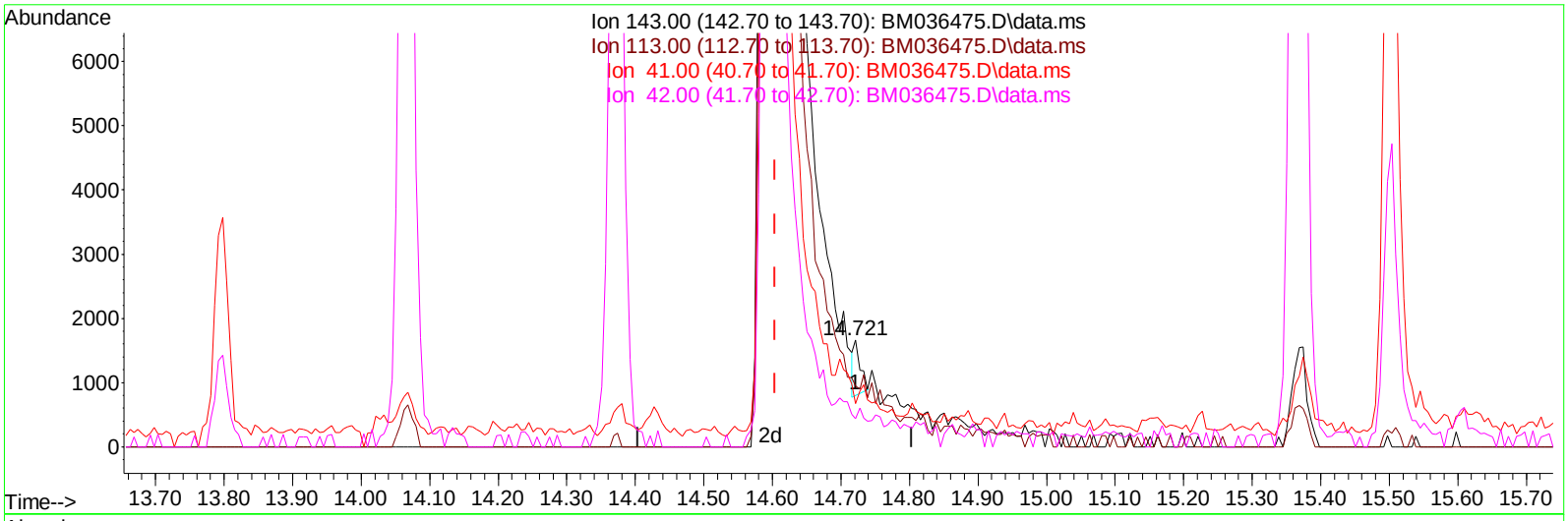
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
Data File : BM036475.D
Acq On : 01 Sep 2022 16:23
Operator : CG/JU
Sample : PB147416BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

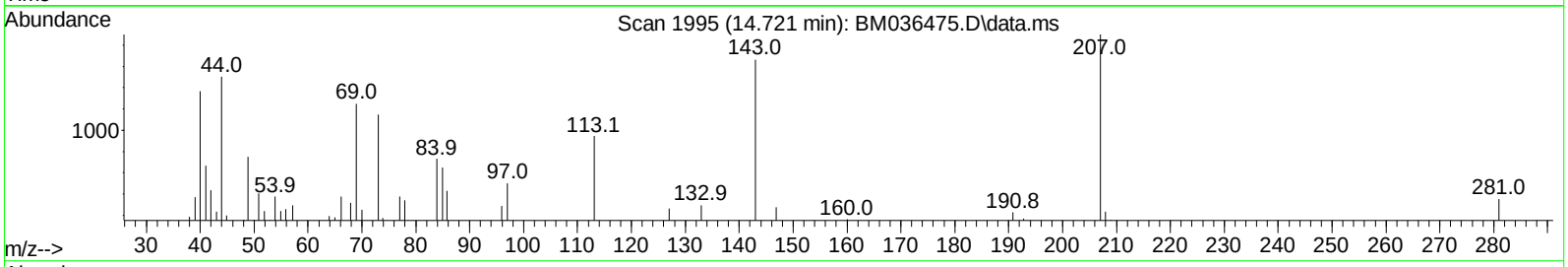
Quant Time: Sep 02 02:21:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Sep 02 01:33:23 2022
Response via : Initial Calibration

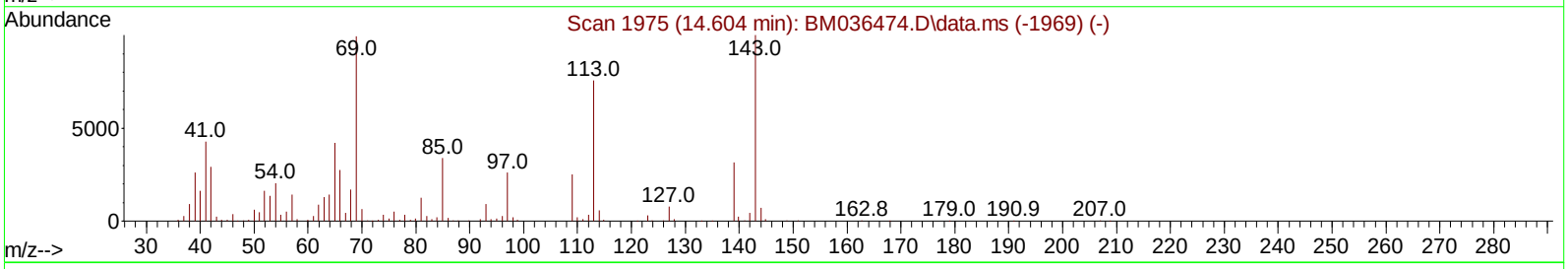
Instrument :
BNA_M
ClientSampleId :
SBLK416

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/02/2022
Supervised By :Sohil Jodhani 09/02/2022







TIC: BM036475.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.721min (+ 0.117) 0.09 ng/ul

response 570

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	56.76
41.00	41.00	40.29
42.00	29.30	26.16

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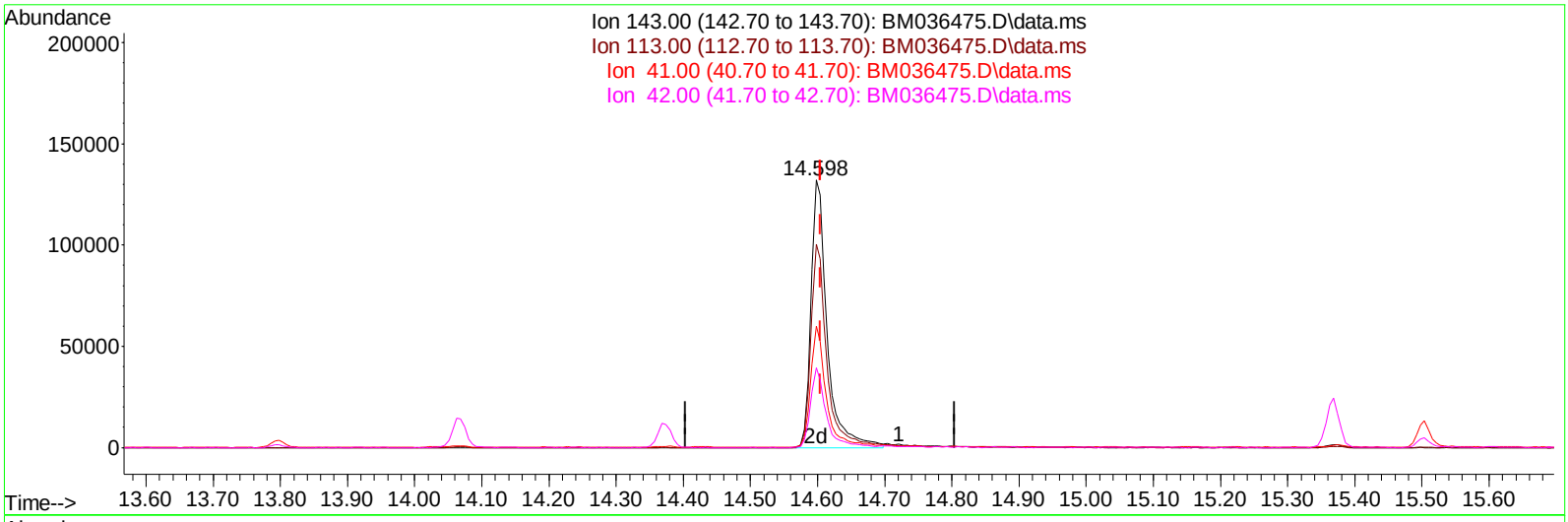

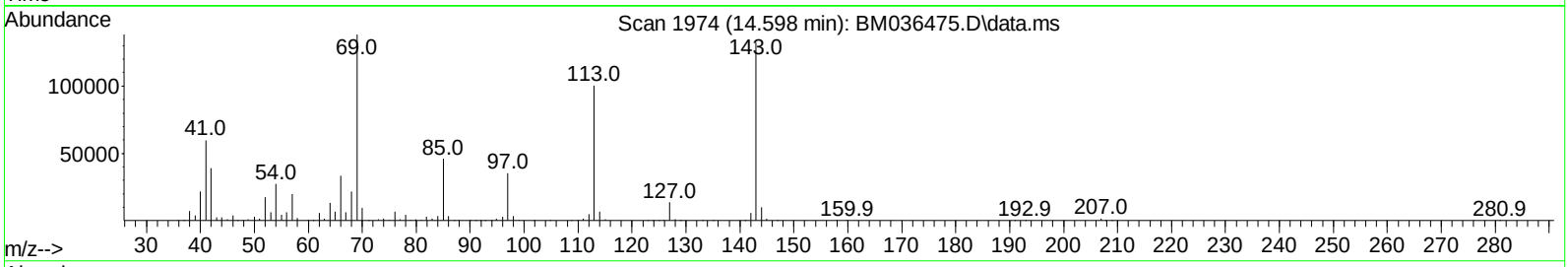

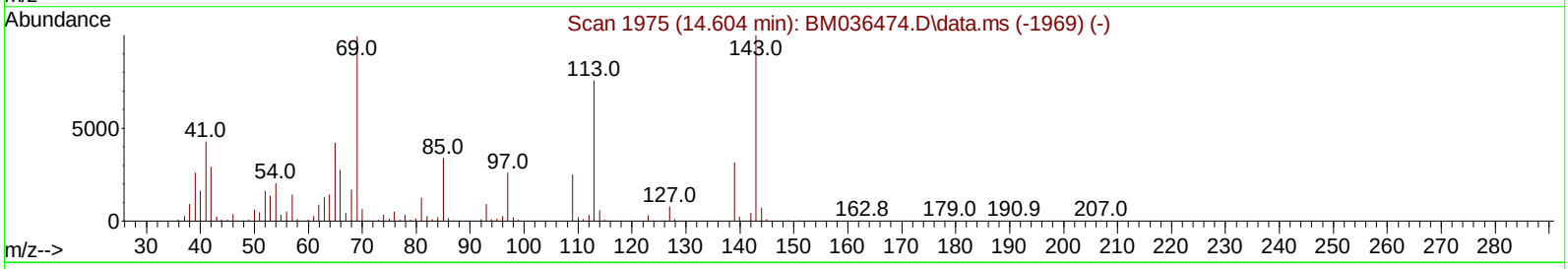

TIC: BM036475.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (-0.006) 34.61 ng/ul m

response 219325

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	76.16#
41.00	41.00	45.30
42.00	29.30	29.73

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	187397	20.000	ng/ul	0.00
20) Naphthalene-d8	10.522	136	853090	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.374	164	528627	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.121	188	1099231	20.000	ng/ul	0.00
79) Chrysene-d12	21.309	240	1112552	20.000	ng/ul	0.00
88) Perylene-d12	23.609	264	1132543	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	30385	6.179	ng/uL	0.00
4) Pyridine-d5	3.551	84	395446	29.163	ng/ul	0.00
7) Phenol-d5	6.898	99	503605	31.483	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	357462	35.502	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.251	132	438195	34.866	ng/ul	0.00
15) 4-Methylphenol-d8	8.445	113	422438	35.370	ng/ul	0.00
21) Nitrobenzene-d5	8.892	128	218167	33.730	ng/ul	0.00
24) 2-Nitrophenol-d4	9.610	143	222641	33.935	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.145	165	379414	32.365	ng/ul	0.00
31) 4-Chloroaniline-d4	10.675	131	341167	18.773	ng/ul	0.00
46) Dimethylphthalate-d6	13.798	166	1306718	39.073	ng/ul	0.00
49) Acenaphthylene-d8	14.068	160	1552411	37.217	ng/ul	0.00
54) 4-Nitrophenol-d4	14.598	143	219325m	34.611	ng/ul	0.00
60) Fluorene-d10	15.368	176	1166761	38.609	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	174876	30.944	ng/ul	0.00
73) Anthracene-d10	17.221	188	1854895	37.950	ng/ul	0.00
81) Pyrene-d10	19.521	212	2196724	38.063	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.462	264	2125807	38.283	ng/ul	0.00
Target Compounds						Qvalue

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

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