Quantitation Report (Qedit)

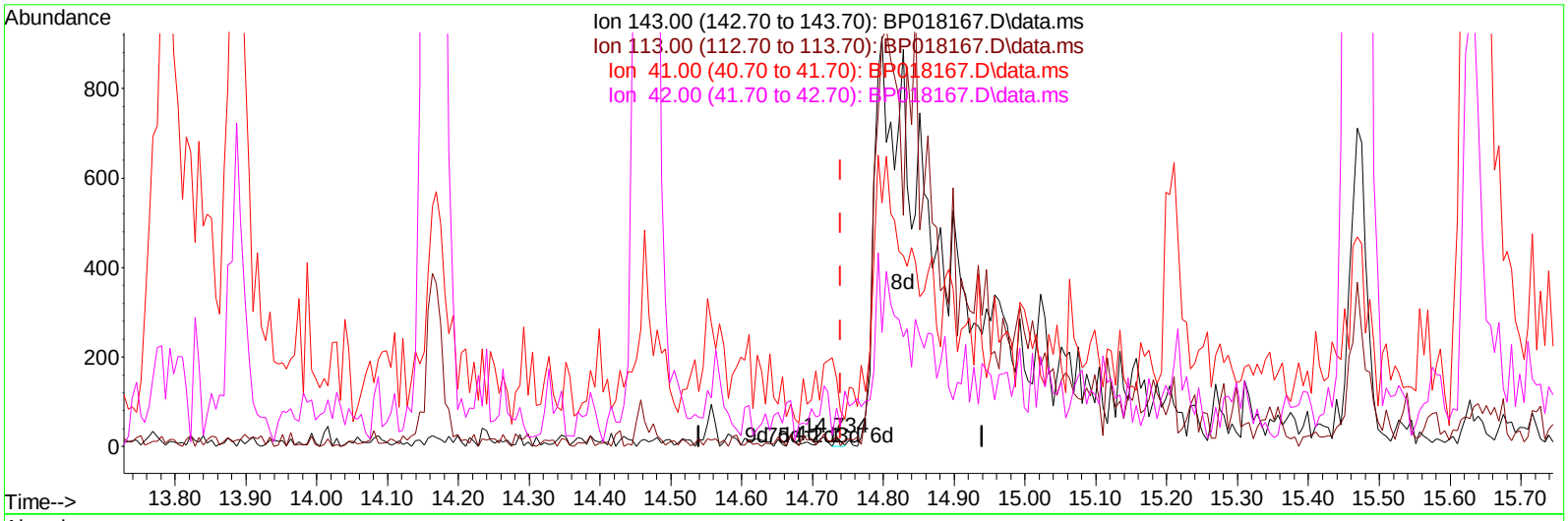
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018167.D
Acq On : 14 Nov 2023 21:47
Operator : MA/JU
Sample : 05337-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

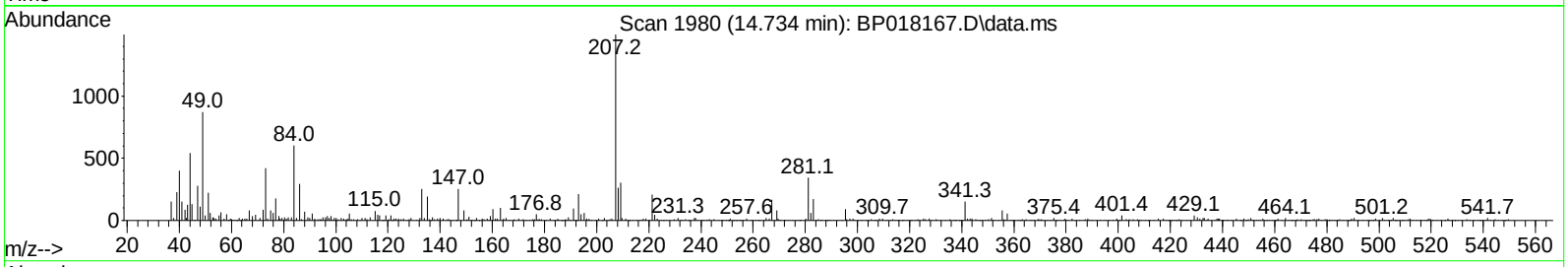
Quant Time: Nov 15 00:40:52 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

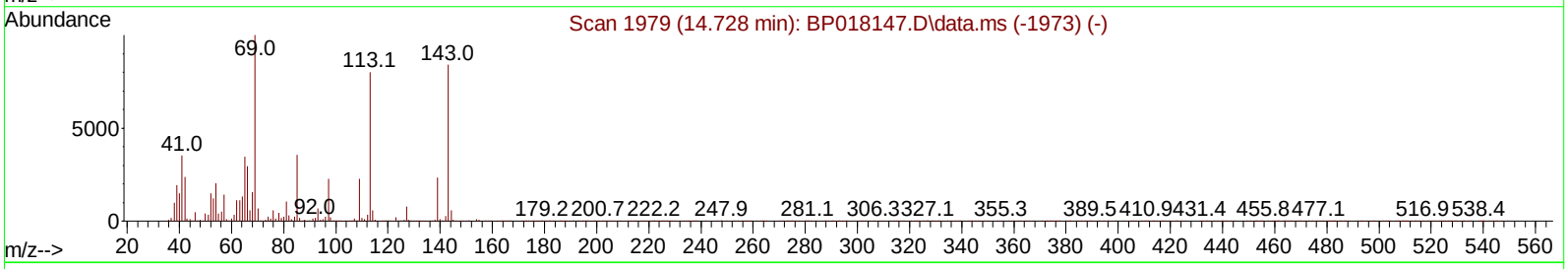
Instrument :
BNA_P
ClientSampleId :
GCDJ8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023







TIC: BP018167.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 0.01 ng/ul

response 13

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	109.52
41.00	47.70	723.81#
42.00	30.10	495.24#

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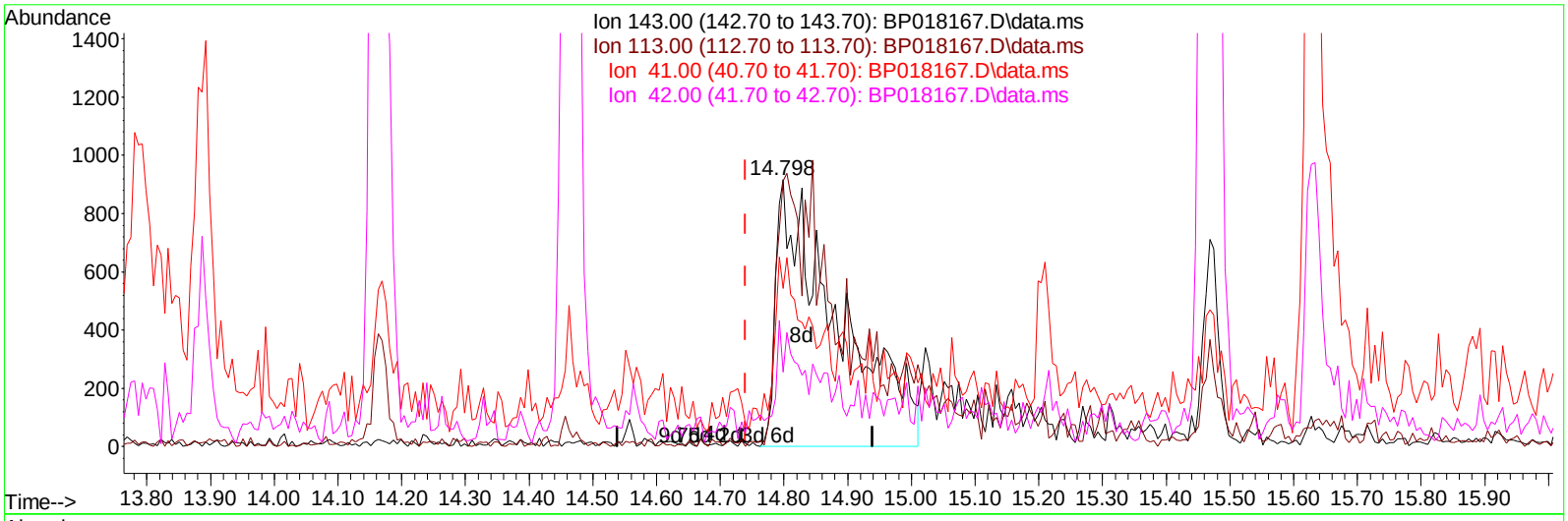

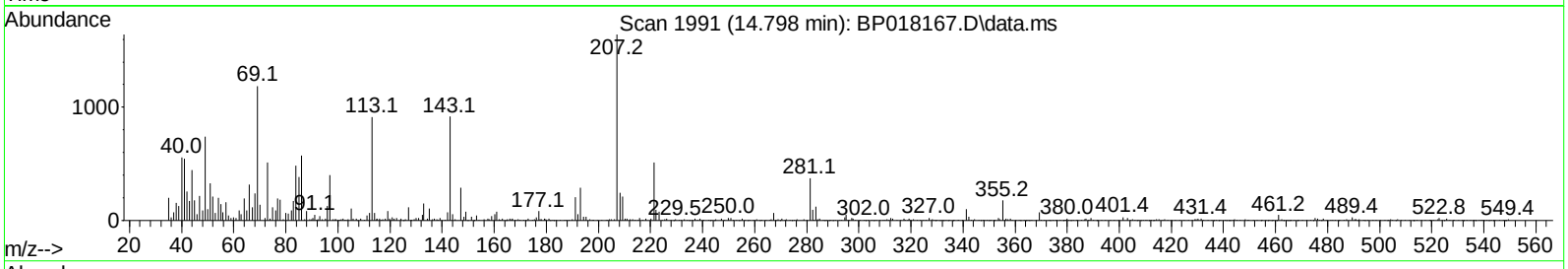

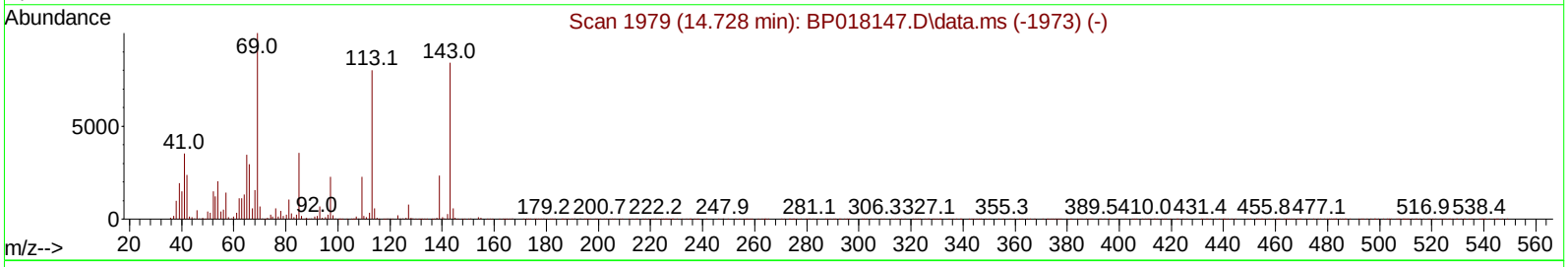

TIC: BP018167.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 2.36 ng/ul m

response 6097

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	99.56
41.00	47.70	59.39#
42.00	30.10	27.84

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	73061	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	295928	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	187525	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.245	188	427661	20.000	ng/ul	-0.01
79) Chrysene-d12	21.380	240	415508	20.000	ng/ul	-0.01
88) Perylene-d12	23.845	264	468933	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	7021	3.390	ng/uL	0.00
4) Pyridine-d5	3.622	84	13379	2.478	ng/ul	0.02
7) Phenol-d5	6.969	99	22889	3.245	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	100834	24.326	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	90845	16.264	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	51965	8.991	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	60716	22.773	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	66262	21.937	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	104660	19.441	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	164686	22.307	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	487112	28.153	ng/ul	-0.01
49) Acenaphthylene-d8	14.163	160	441774	23.442	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	6097m	2.358	ng/ul	0.06
60) Fluorene-d10	15.469	176	391195	27.385	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	58774	19.010	ng/ul	-0.03
73) Anthracene-d10	17.345	188	686665	29.688	ng/ul	-0.01
81) Pyrene-d10	19.604	212	874673	31.515	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	875239	32.099	ng/ul	-0.02
Target Compounds						Qvalue

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

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