Quantitation Report (Qedit)

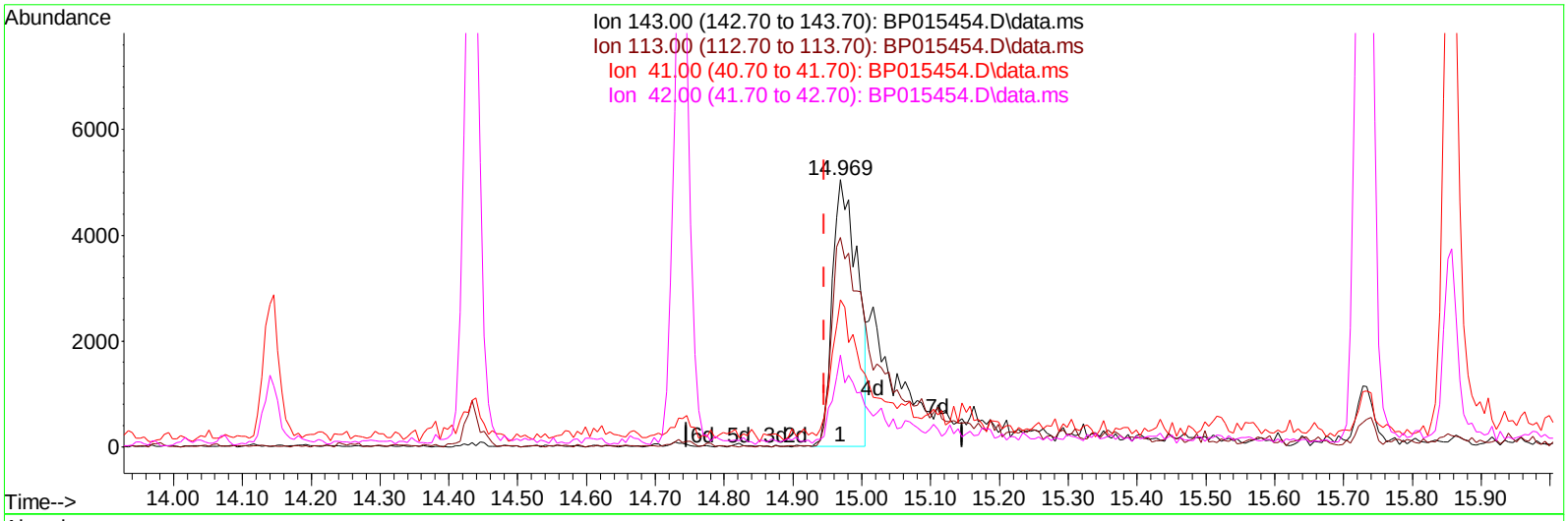
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
Data File : BP015454.D
Acq On : 09 Jun 2023 16:19
Operator : MA/JU
Sample : 03097-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

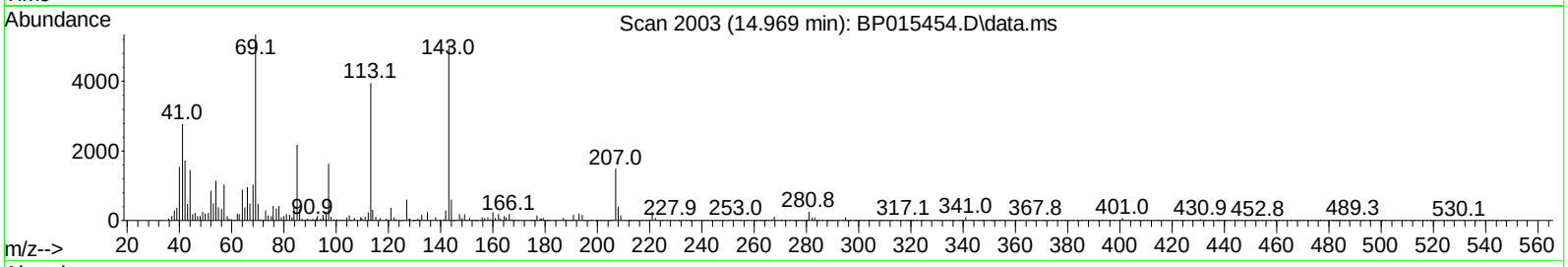
Quant Time: Jun 09 22:58:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 07 23:46:15 2023
Response via : Initial Calibration

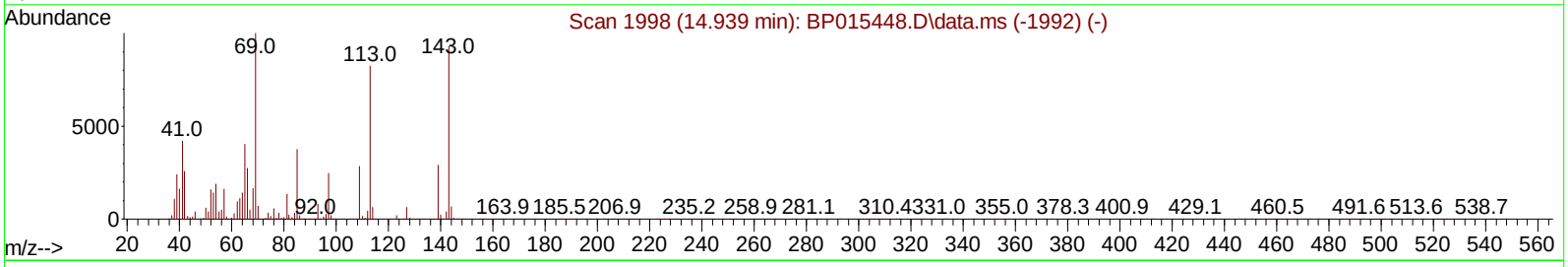
Instrument :
BNA_P
ClientSampleId :
EZEN6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/12/2023
Supervised By :mohammad ahmed 06/13/2023







TIC: BP015454.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.024) 2.10 ng/ul

response 12740

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	78.37
41.00	42.20	54.96#
42.00	29.40	34.22

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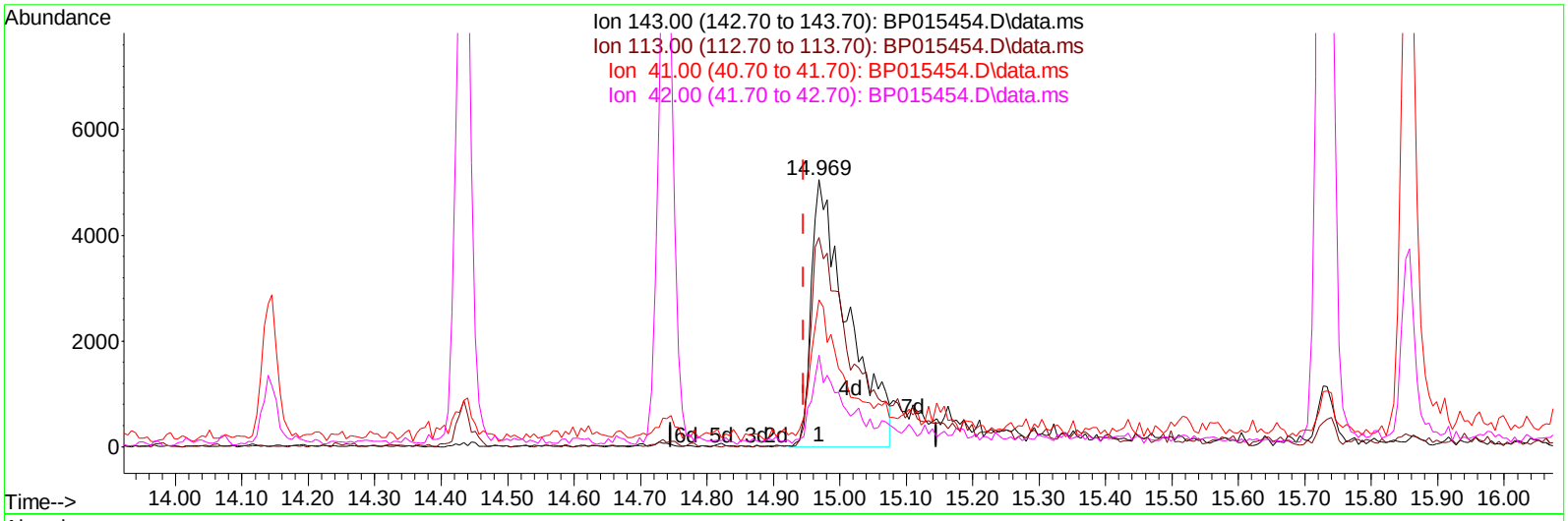

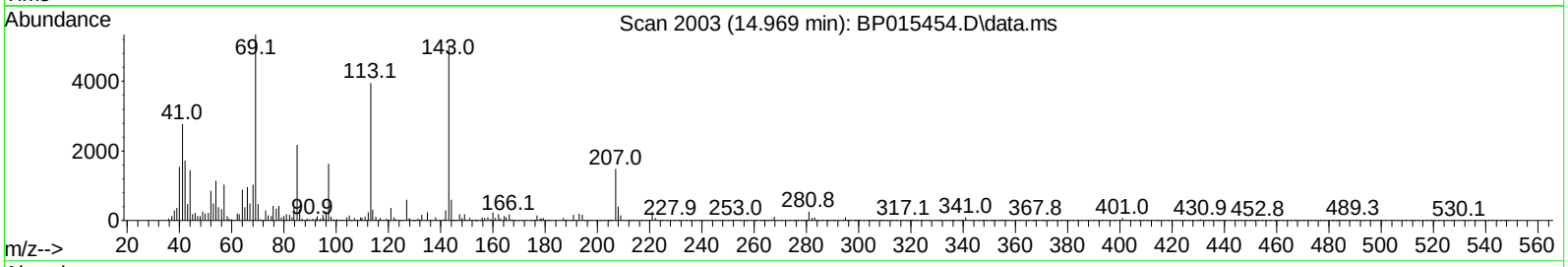

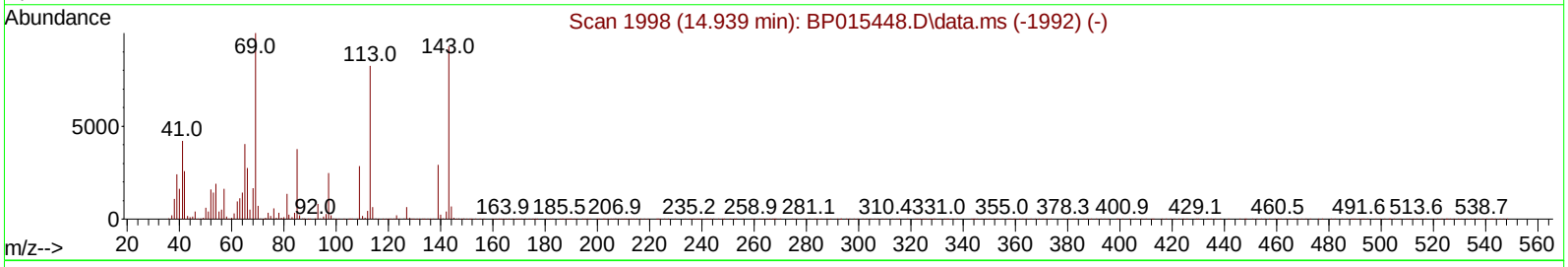

TIC: BP015454.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.024) 3.18 ng/ul m

response 19307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	78.37
41.00	42.20	54.96#
42.00	29.40	34.22

Quantitation Report (QT Reviewed)

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Acq On : 09 Jun 2023 16:19
Operator : MA/JU
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Quant Time: Jun 09 23:16:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	134013	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.922	136	619623	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.740	164	429493	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	1011348	20.000	ng/ul	-0.01
79) Chrysene-d12	21.575	240	1023077	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	1091394	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	6353	1.755	ng/uL	0.00
4) Pyridine-d5	3.858	84	81868	8.709	ng/ul	0.00
7) Phenol-d5	7.252	99	63037	5.106	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	220620	29.921	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	199470	22.284	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	140968	13.912	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	141657	29.120	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	155494	27.990	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	262235	25.877	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.063	131	342432	23.541	ng/ul	0.00
46) Dimethylphthalate-d6	14.146	166	1156691	34.153	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	1194339	32.250	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	19307m	3.183	ng/ul	0.02
60) Fluorene-d10	15.734	176	962719	33.709	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	189783	29.621	ng/ul	0.00
73) Anthracene-d10	17.592	188	1677481	36.455	ng/ul	0.00
81) Pyrene-d10	19.816	212	2114197	38.611	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	2061203	37.640	ng/ul	-0.01

Target Compounds Qvalue

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

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