Quantitation Report (Qedit)

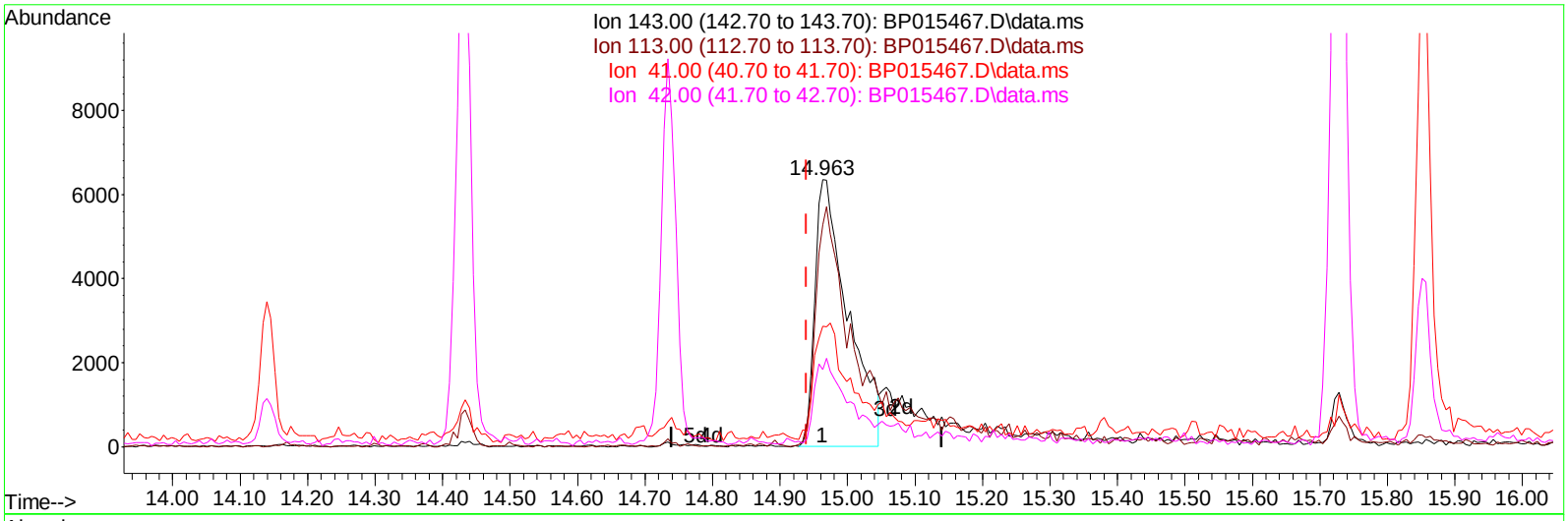
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
Data File : BP015467.D
Acq On : 10 Jun 2023 00:44
Operator : MA/JU
Sample : 03097-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

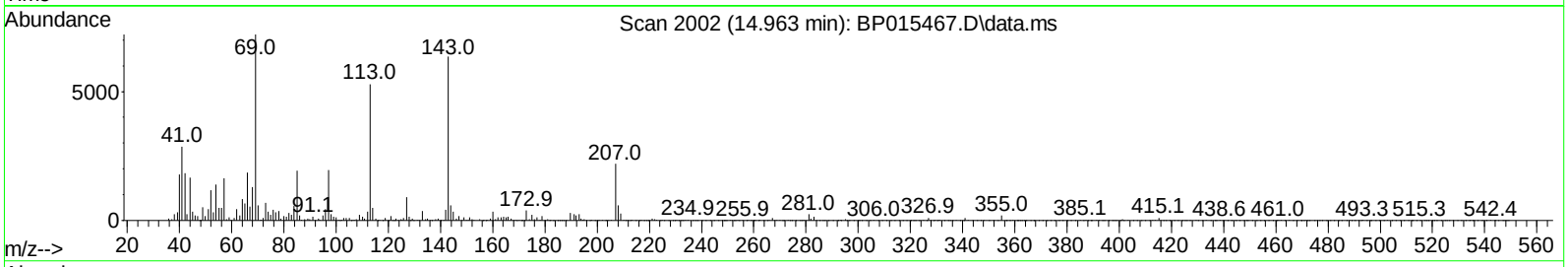
Quant Time: Jun 10 01:22:20 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

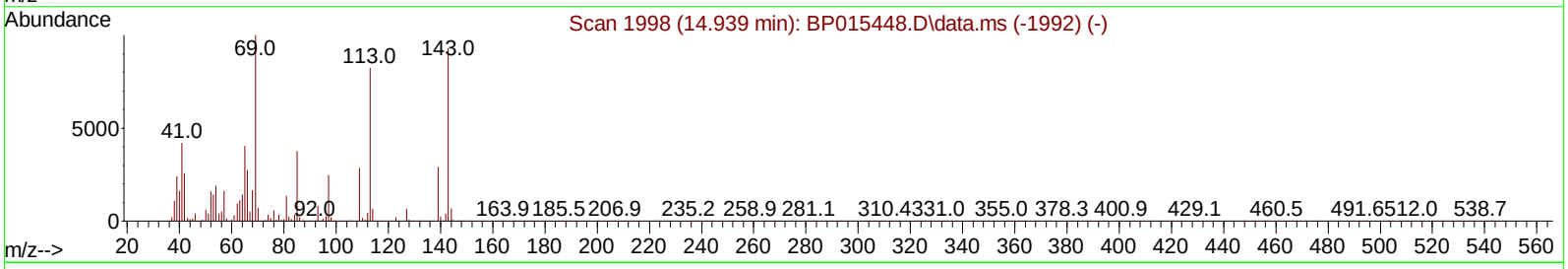
Instrument :
BNA_P
ClientSampleId :
EZEP1

Manual IntegrationsAPPROVED

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







TIC: BP015467.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 4.16 ng/ul

response 21755

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	83.17
41.00	42.20	45.23
42.00	29.40	28.87

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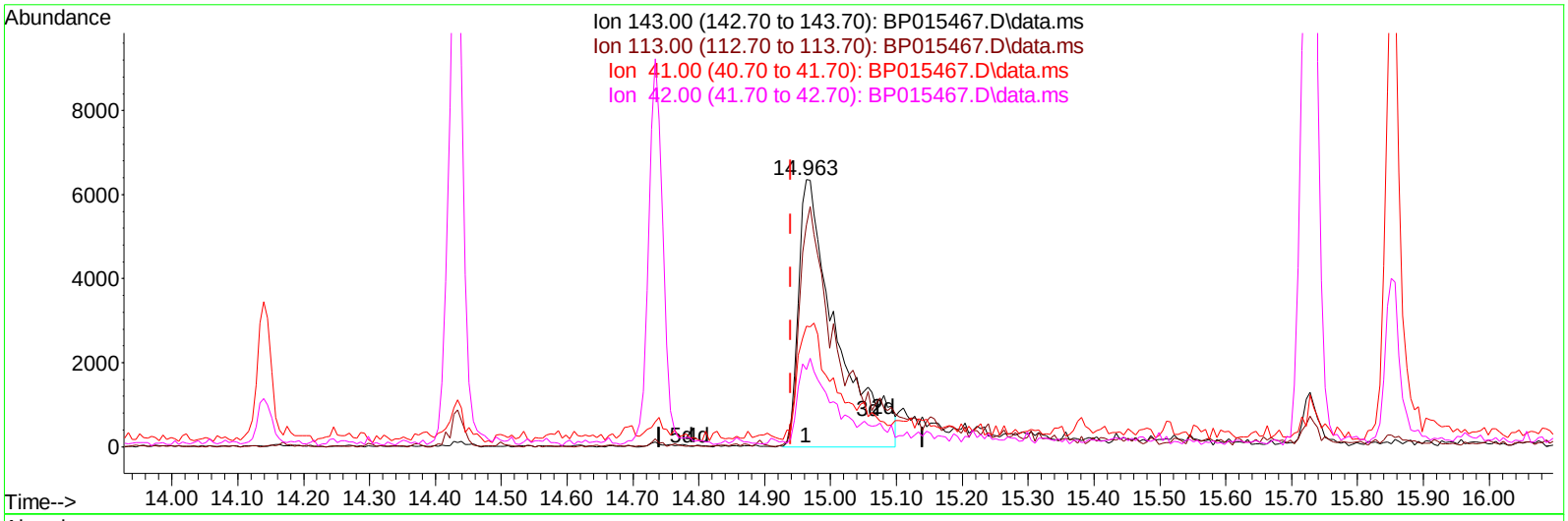

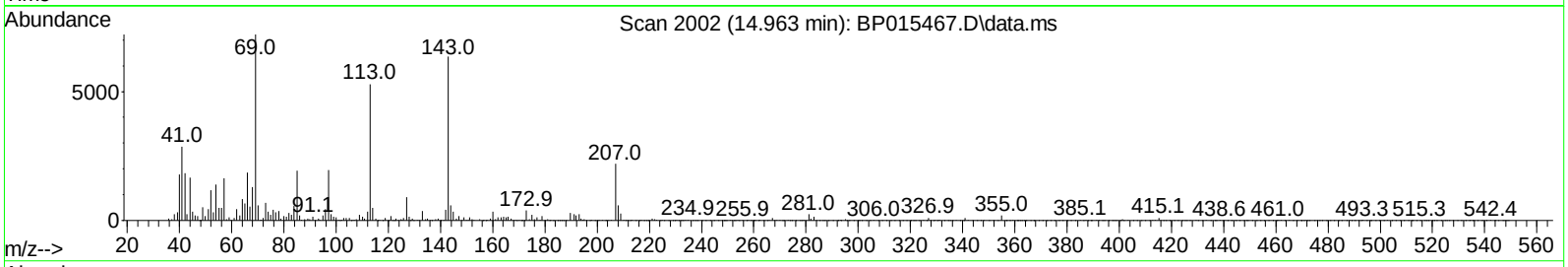

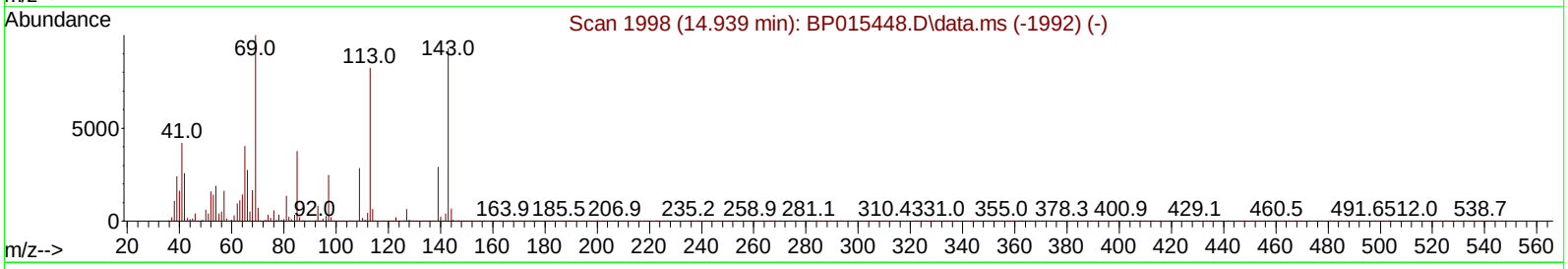

TIC: BP015467.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 4.82 ng/ul m

response 25241

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	83.17
41.00	42.20	45.23
42.00	29.40	28.87

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	112568	20.000	ng/ul	0.00
20) Naphthalene-d8	10.922	136	518137	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	370712	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	876543	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	919147	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	964691	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	7950	2.615	ng/uL	0.00
4) Pyridine-d5	3.864	84	51261	6.492	ng/ul	0.02
7) Phenol-d5	7.252	99	75788	7.308	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	237350	38.323	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	221308	29.434	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	156076	18.338	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	157653	38.756	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	172298	37.089	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	280337	33.081	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	368819	30.321	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	1218858	41.695	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	1301749	40.724	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143	25241m	4.822	ng/ul	0.02
60) Fluorene-d10	15.728	176	1034246	41.955	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	194590	35.043	ng/ul	0.00
73) Anthracene-d10	17.592	188	1818140	45.589	ng/ul	0.00
81) Pyrene-d10	19.816	212	2255106	45.841	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	2248782	46.458	ng/ul	0.00

Target Compounds Qvalue

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

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