Quantitation Report (Qedit)

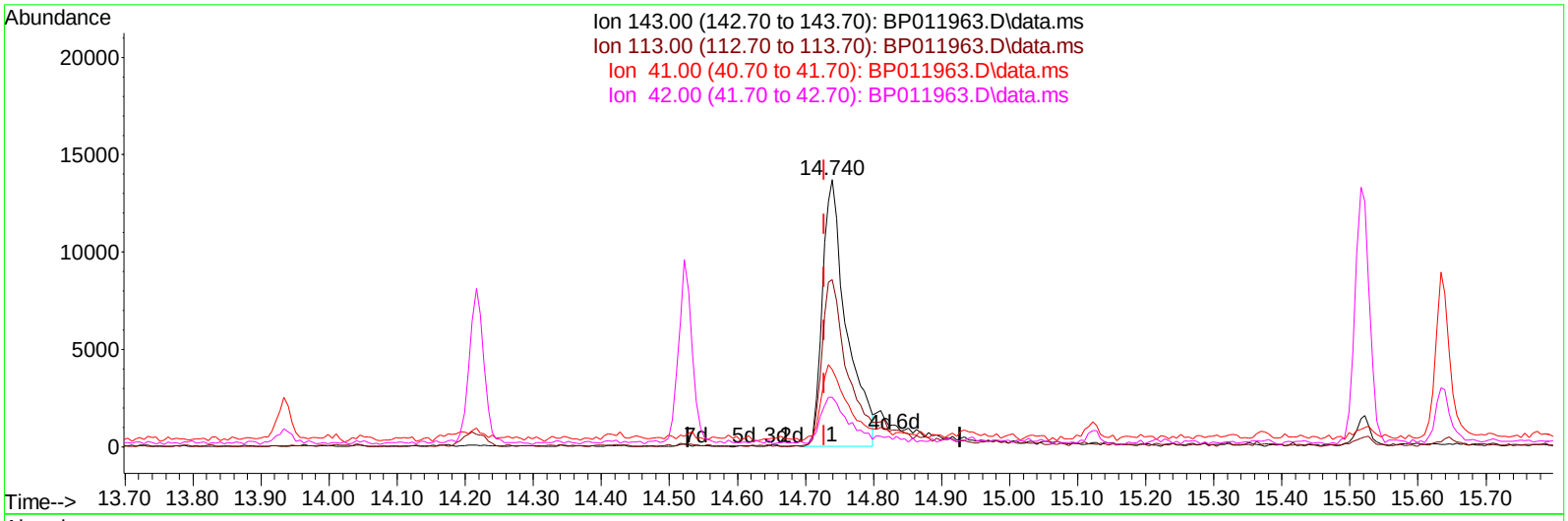
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011963.D
Acq On : 06 Oct 2022 11:03
Operator : CG/JU
Sample : N4956-11
Misc :
ALS Vial : 44 Sample Multiplier: 1

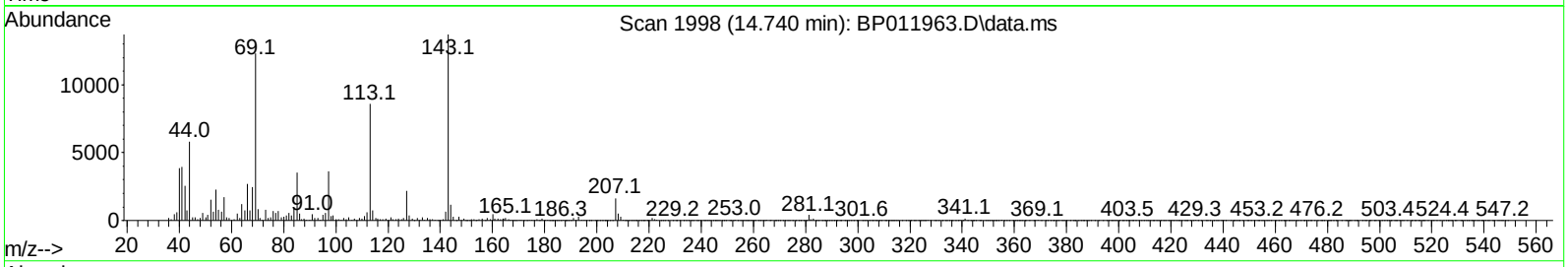
Quant Time: Oct 06 23:04:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 05 23:18:55 2022
Response via : Initial Calibration

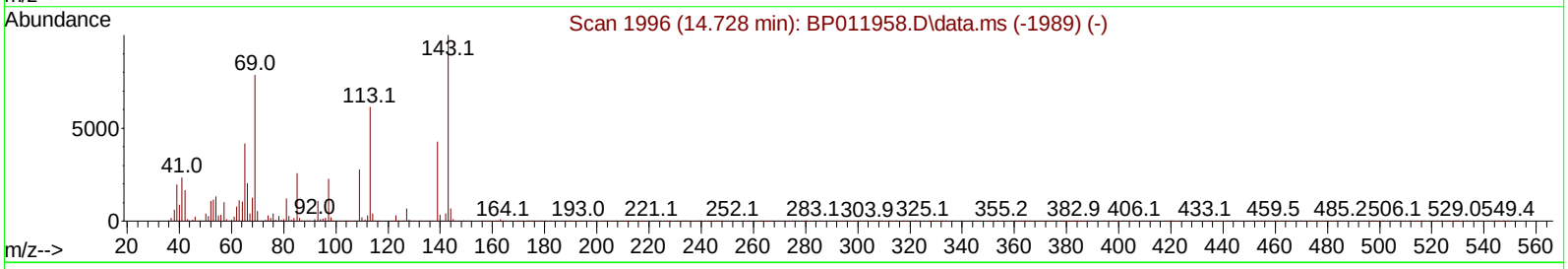
Instrument :
BNA_P
ClientSampleId :
CB8T1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/07/2022
Supervised By :mohammad ahmed 10/10/2022







TIC: BP011963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.37 ng/ul

response 33573

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	62.71
41.00	24.20	28.85
42.00	16.40	18.55

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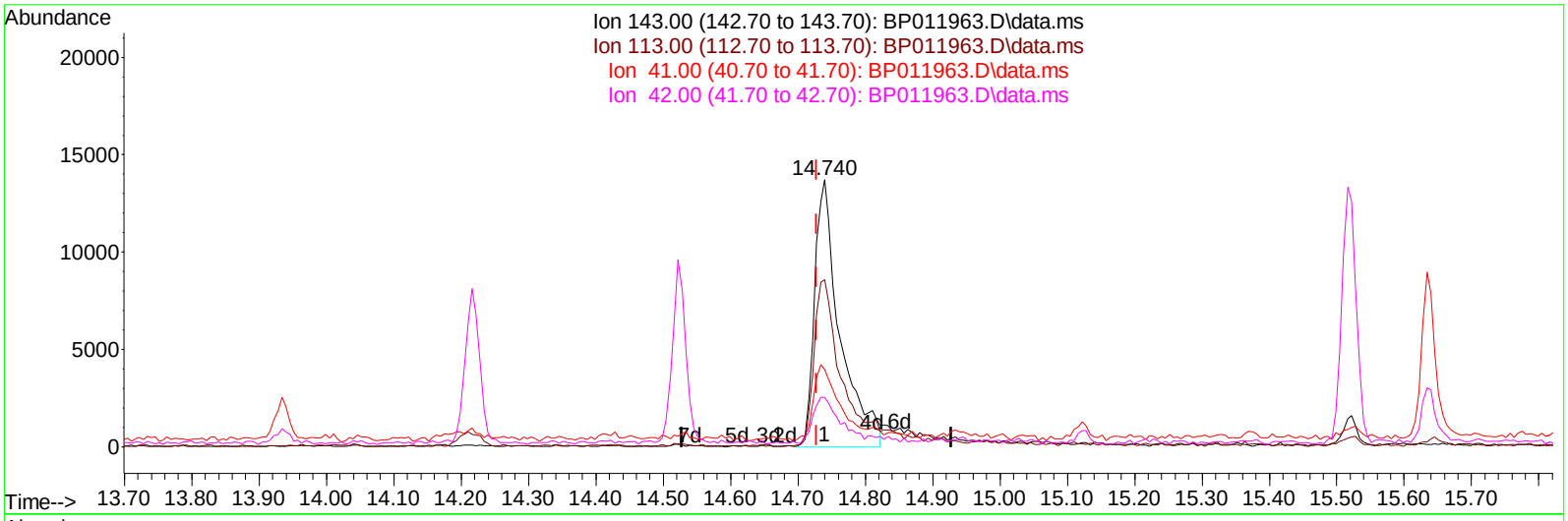

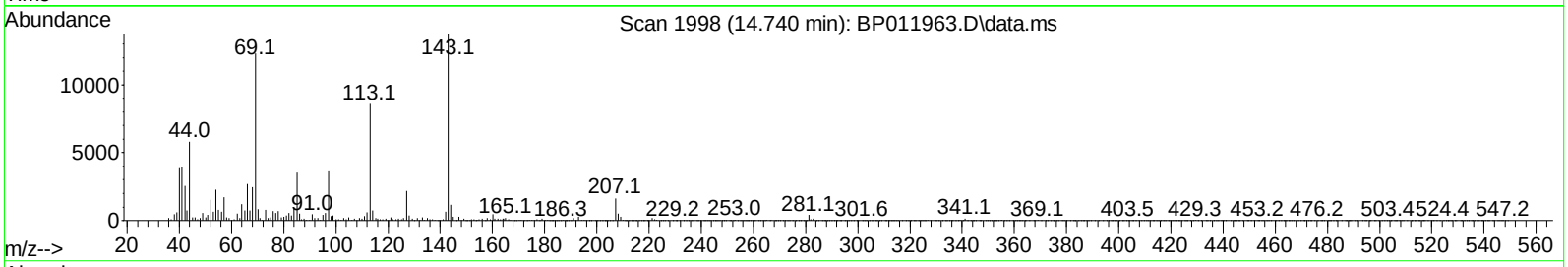

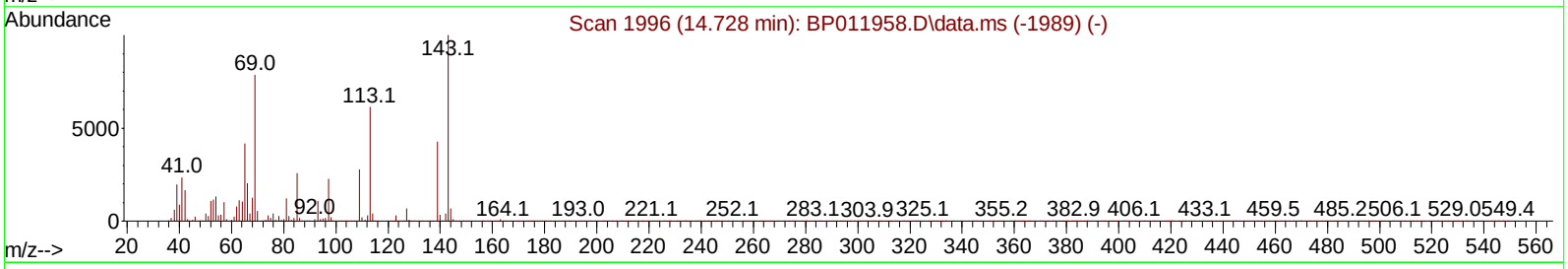

TIC: BP011963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.60 ng/ul m

response 35829

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	62.71
41.00	24.20	28.85
42.00	16.40	18.55

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	261606	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1107075	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	672261	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1401448	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1331257	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1333286	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	22634	3.816	ng/uL	0.00
4) Pyridine-d5	3.728	84	91400	5.260	ng/ul	0.00
7) Phenol-d5	7.040	99	108157	4.789	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	303855	24.273	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	335062	18.794	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	207339	11.106	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	223188	26.468	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	214488	23.774	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	368240	22.067	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	323687	12.724	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1335299	27.776	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1484096	27.080	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	35829m	3.599	ng/ul	0.01
60) Fluorene-d10	15.522	176	1191308	28.535	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	174875	20.616	ng/ul	0.00
73) Anthracene-d10	17.381	188	1997491	31.290	ng/ul	0.00
81) Pyrene-d10	19.616	212	2330667	34.569	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2205464	33.044	ng/ul	0.00
Target Compounds						Qvalue

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

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