Quantitation Report (Qedit)

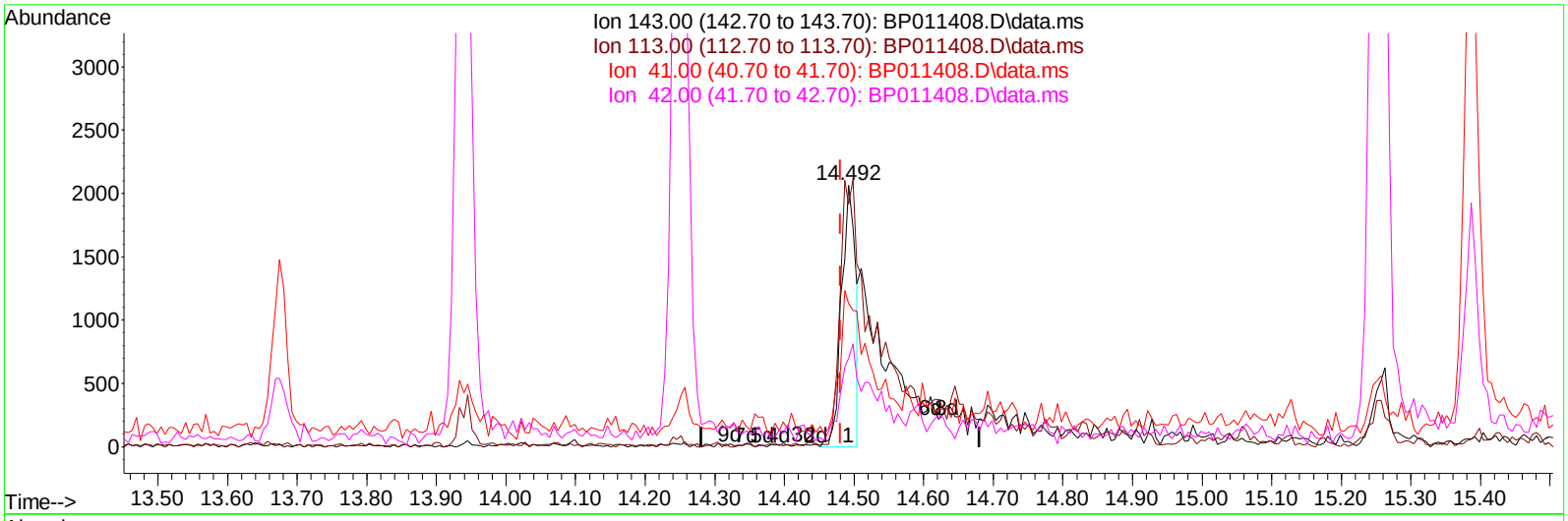
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011408.D
Acq On : 12 Aug 2022 03:42
Operator : CG/JU
Sample : N4128-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

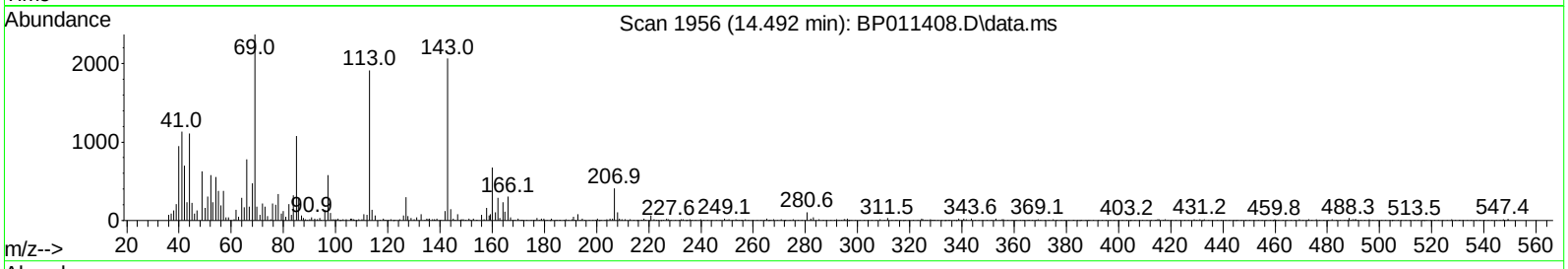
Quant Time: Aug 12 04:08:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

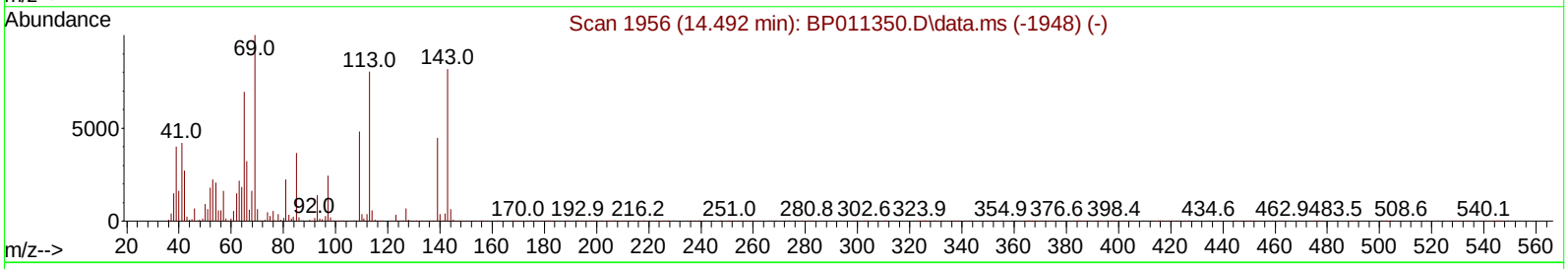
Instrument :
BNA_P
ClientSampleId :
BGBT4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011408.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.012) 1.32 ng/ul

response 2922

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	92.60#
41.00	54.40	54.72
42.00	35.50	33.82

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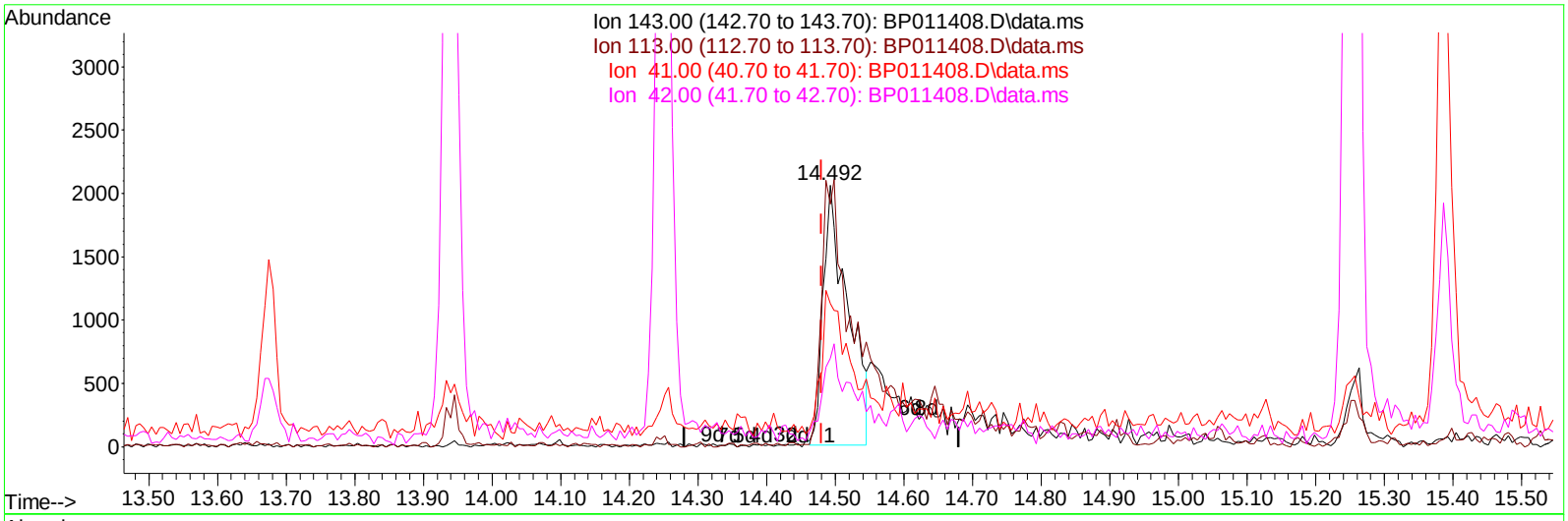

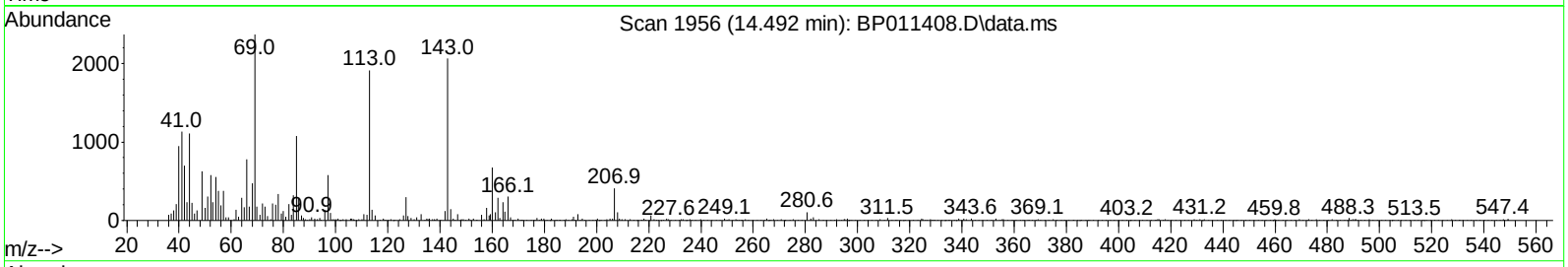

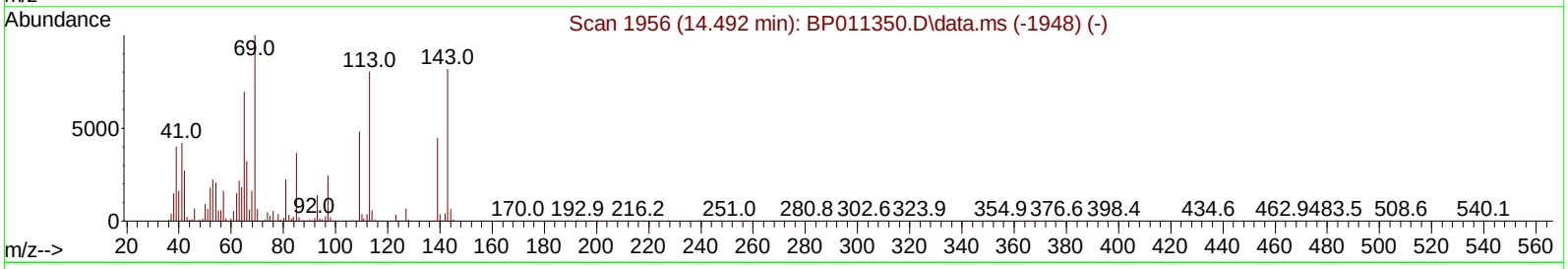

TIC: BP011408.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.012) 2.34 ng/ul m

response 5163

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	92.60#
41.00	54.40	54.72
42.00	35.50	33.82

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	72878	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	308893	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	174335	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	363324	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	356041	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	378102	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	9777	4.113	ng/uL	0.00
4) Pyridine-d5	3.499	84	37841	6.310	ng/ul	0.01
7) Phenol-d5	6.763	99	32944	5.108	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	123057	26.909	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	107927	21.328	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	63039	12.448	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	69252	32.942	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	64469	33.496	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	95795	24.646	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	134278	19.761	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	397910	32.895	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	459446	31.553	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	5163m	2.337	ng/ul	0.01
60) Fluorene-d10	15.257	176	328465	31.752	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	47658	28.240	ng/ul	0.00
73) Anthracene-d10	17.122	188	597645	36.597	ng/ul	0.00
81) Pyrene-d10	19.386	212	706912	36.038	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	710568	37.265	ng/ul	0.00

Target Compounds Qvalue

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

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