Quantitation Report (Qedit)

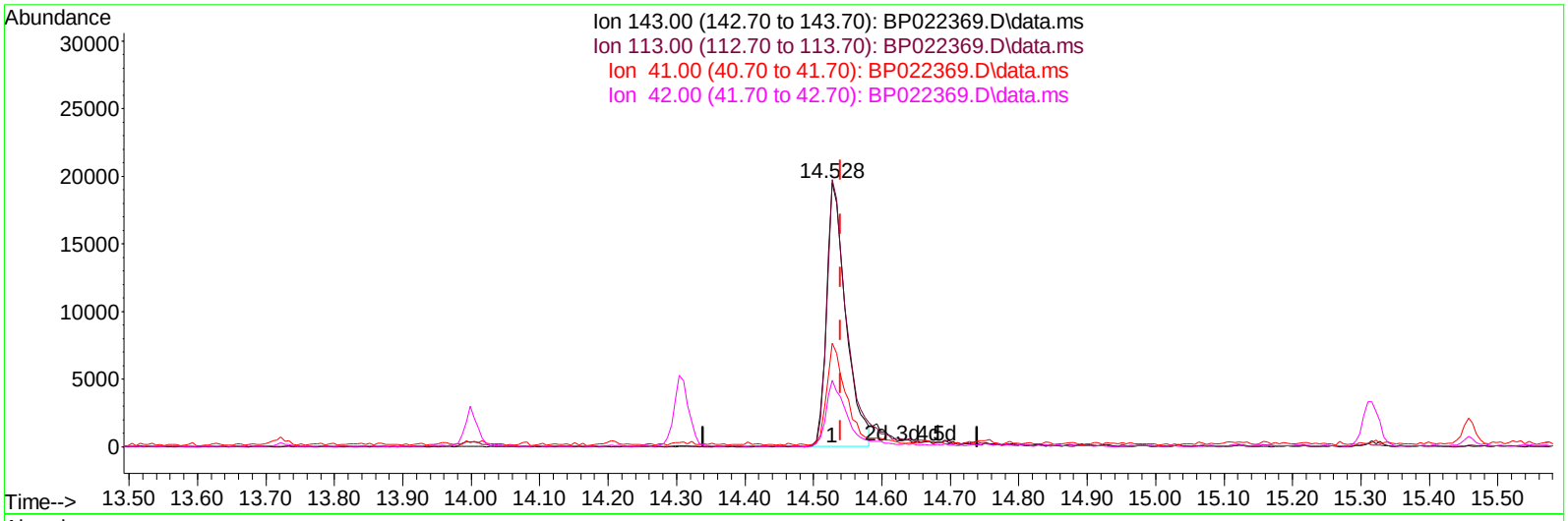
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022369.D
Acq On : 15 Oct 2024 00:10
Operator : RC/JU
Sample : P4237-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

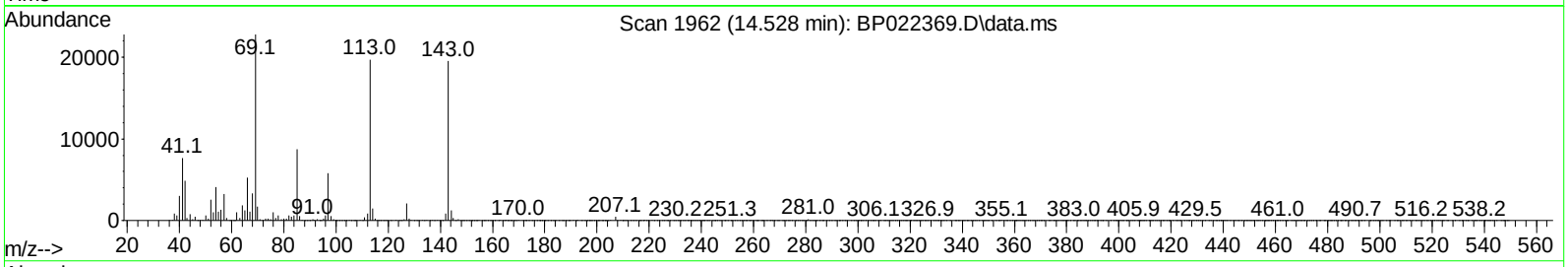
Quant Time: Oct 15 01:54:07 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 14 11:01:55 2024
Response via : Initial Calibration

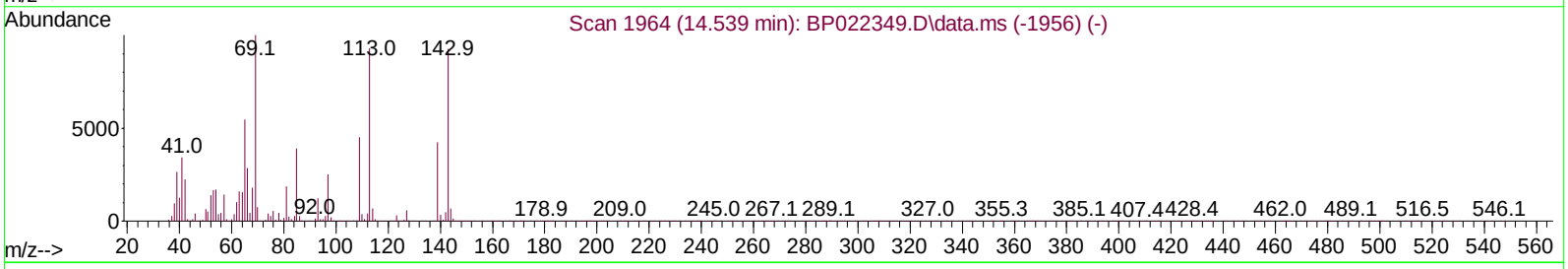
Instrument :
BNA_P
ClientSampleId :
A4F11

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/15/2024
Supervised By :mohammad ahmed 10/22/2024







TIC: BP022369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (-0.012) 9.12 ng/ul

response 38217

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	101.07
41.00	35.90	39.31
42.00	23.70	25.07

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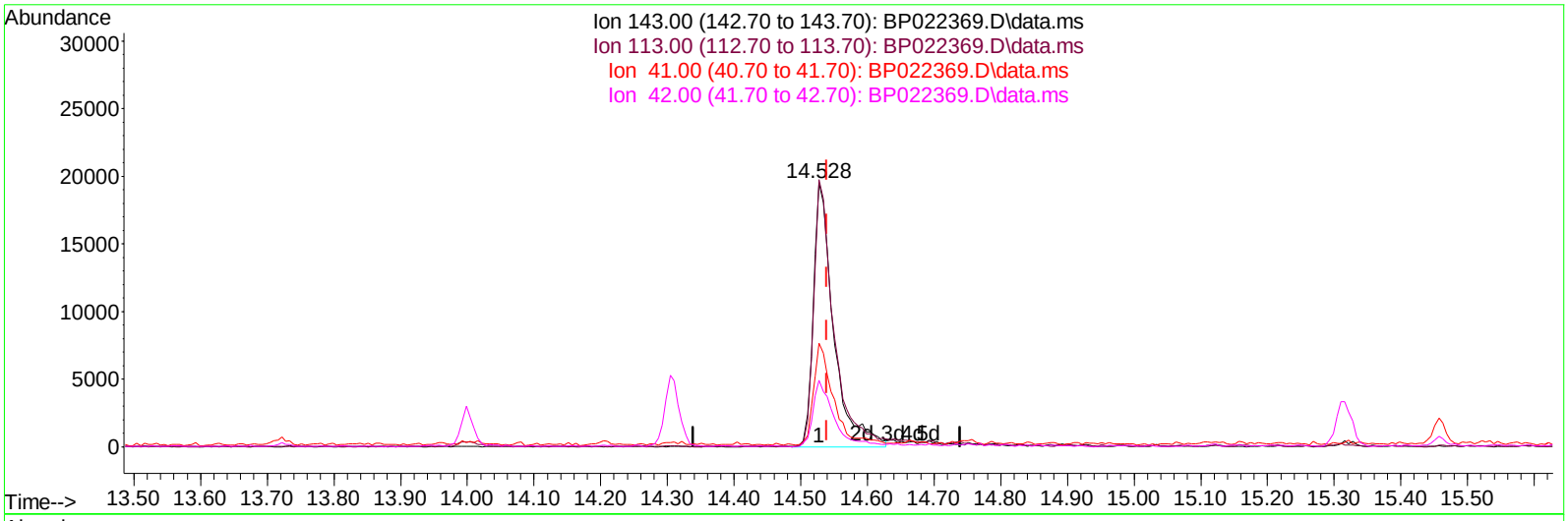

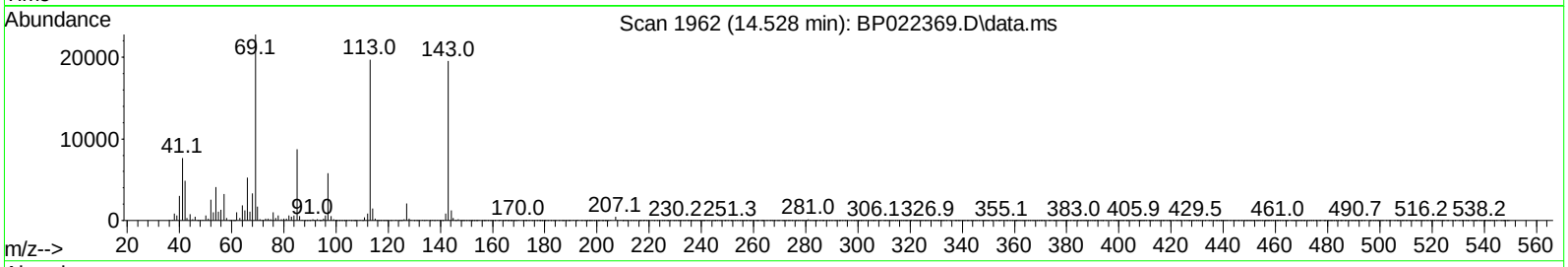

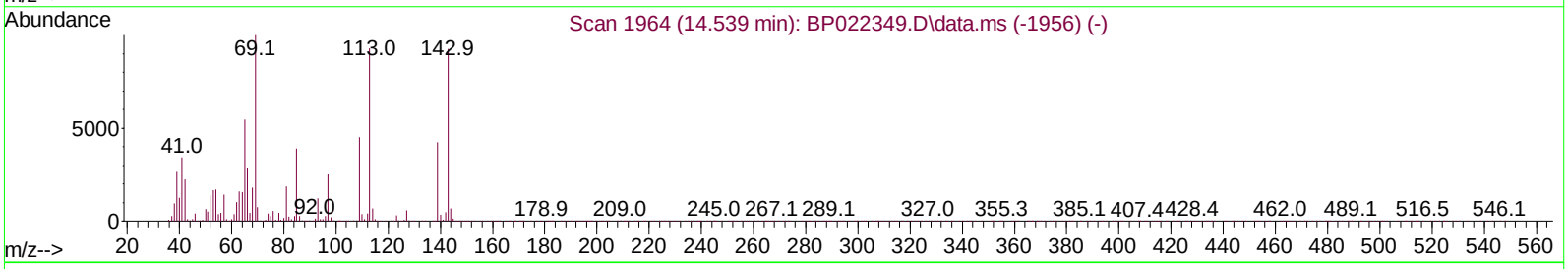

TIC: BP022369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (-0.012) 9.81 ng/ul m

response 41101

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	101.07
41.00	35.90	39.31
42.00	23.70	25.07

Quantitation Report (QT Reviewed)

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Acq On : 15 Oct 2024 00:10
Operator : RC/JU
Sample : P4237-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 15 01:57:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 14 11:01:55 2024
Response via : Initial Calibration

Instrument :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	102872	20.000	ng/ul	0.00
20) Naphthalene-d8	10.451	136	395868	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.304	164	314310	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	779885	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	807519	20.000	ng/ul	-0.01
88) Perylene-d12	24.845	264	941055	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	5016	1.895	ng/uL	0.00
4) Pyridine-d5	3.646	84	8941	1.230	ng/ul	0.01
7) Phenol-d5	6.887	99	97000	11.202	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	53680	10.547	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	73508	11.963	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	73716	10.667	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	34597	11.366	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	42027	11.926	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	89395	11.936	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	88258	9.973	ng/ul	0.01
46) Dimethylphthalate-d6	13.722	166	305844	12.247	ng/ul	0.01
49) Acenaphthylene-d8	14.004	160	326323	12.303	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	41101m	9.811	ng/ul	-0.01
60) Fluorene-d10	15.316	176	286711	13.237	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	37656	7.342	ng/ul	0.01
73) Anthracene-d10	17.210	188	461314	12.574	ng/ul	0.00
81) Pyrene-d10	19.592	212	545170	12.767	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	461748	9.188	ng/ul	-0.04
Target Compounds						Qvalue
16) Acetophenone	8.505	105	12273	1.076	ng/ul	95
80) Fluoranthene	19.245	202	71927	1.465	ng/ul	99
82) Pyrene	19.622	202	58909	1.120	ng/ul	98

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

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