

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021247.D
Acq On : 30 Jul 2024 23:35
Operator : MA/JU
Sample : P3347-11
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AX3

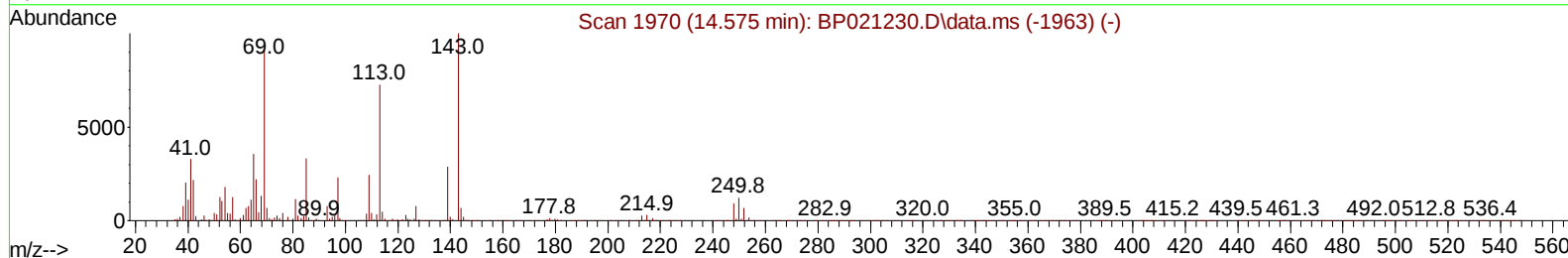
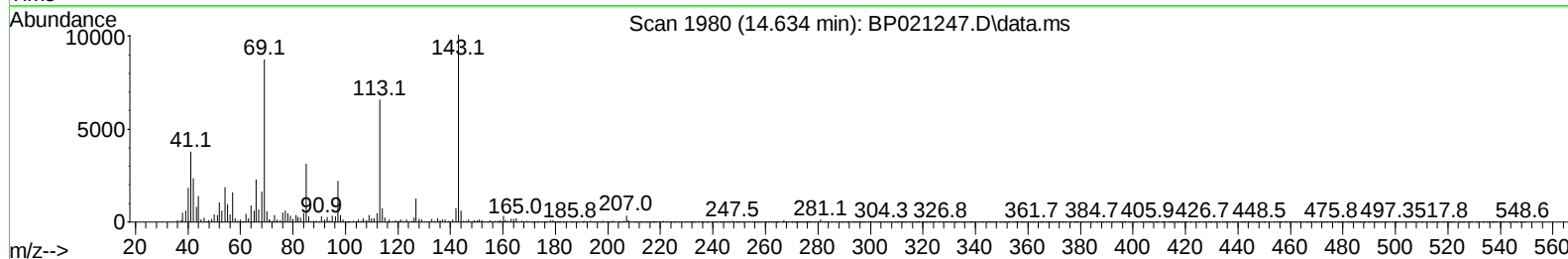
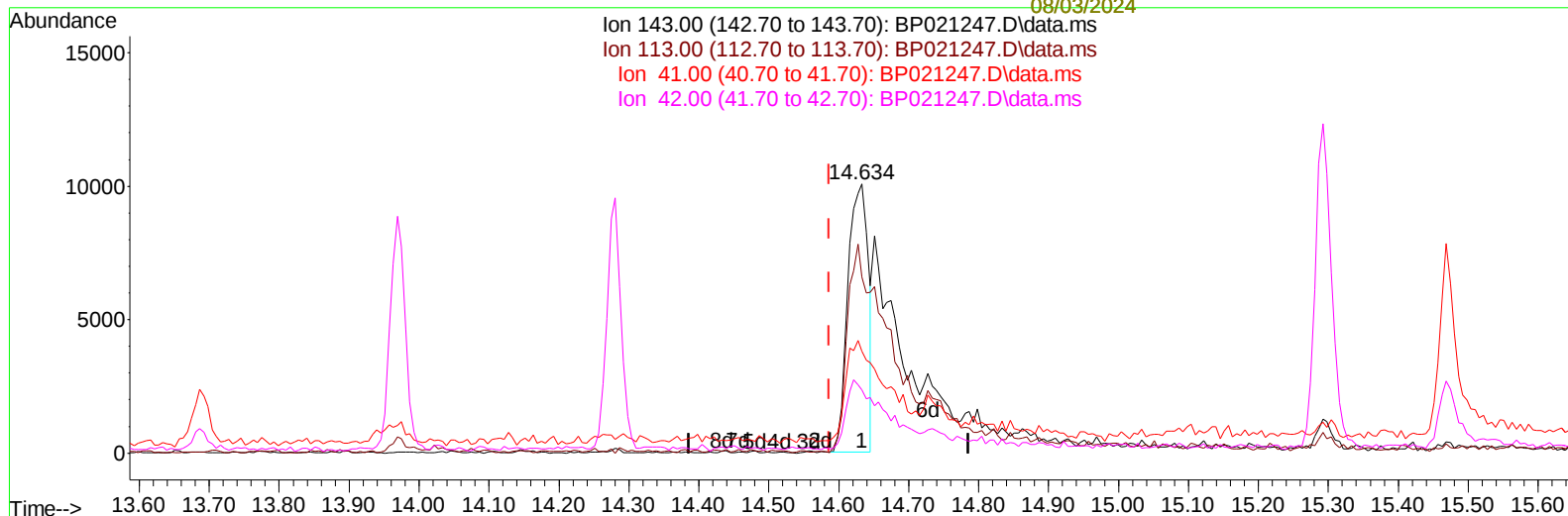
Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/31/2024
Supervised By :mohammad ahmed 09/04/2024

Supervised By :mohammad
ahmed

08/03/2024

Quant Time: Jul 31 00:01:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration



TIC: BP021247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.047) 3.60 ng/u1

response 20612

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	65.35
41.00	37.90	37.80
42.00	27.10	23.53

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	204972	20.000	ng/ul	0.00
20) Naphthalene-d8	10.398	136	822923	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164	553931	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.098	188	1169894	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1017896	20.000	ng/ul	0.01
88) Perylene-d12	24.845	264	1073671	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	18365	4.077	ng/uL	0.00
4) Pyridine-d5	3.599	84	90705	6.956	ng/ul	0.00
7) Phenol-d5	6.863	99	164875	9.866	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67	226456	22.361	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	357795	25.987	ng/ul	0.01
15) 4-Methylphenol-d8	8.369	113	289746	20.876	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	166990	26.126	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	197925	27.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	379891	28.718	ng/ul	0.01
31) 4-Chloroaniline-d4	10.569	131	372491	21.255	ng/ul	0.01
46) Dimethylphthalate-d6	13.687	166	1197309	29.734	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	1299651	28.744	ng/ul	0.00
54) 4-Nitrophenol-d4	14.634	143	49493m	8.639	ng/ul	0.05
60) Fluorene-d10	15.292	176	998175	30.215	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	178289	25.187	ng/ul	0.00
73) Anthracene-d10	17.198	188	1581087	30.428	ng/ul	0.00
81) Pyrene-d10	19.586	212	1879044	29.859	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.627	264	1655489	31.263	ng/ul	0.03

Target Compounds Qvalue

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

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