

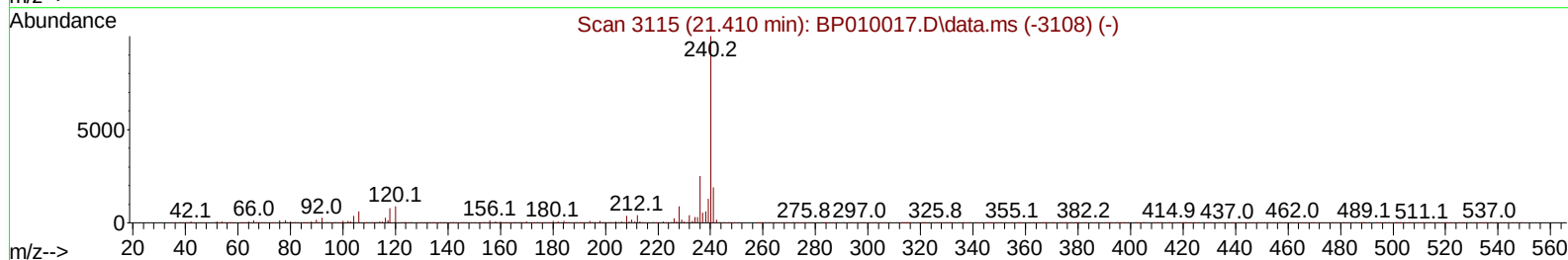
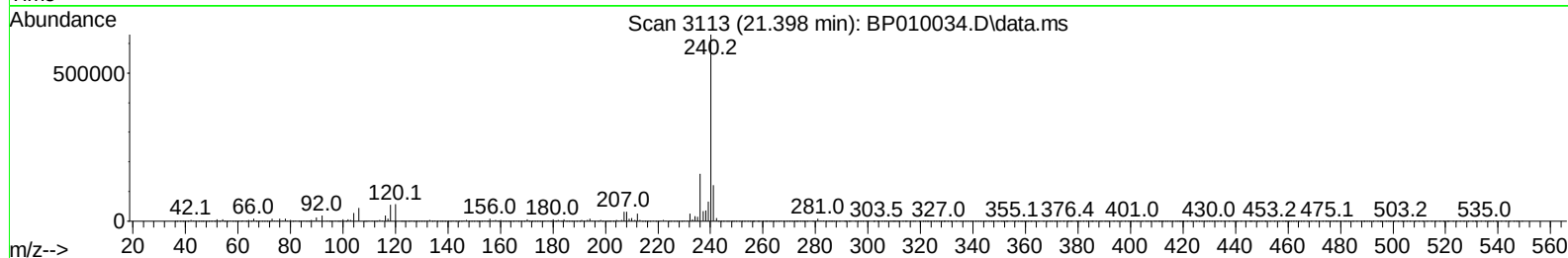
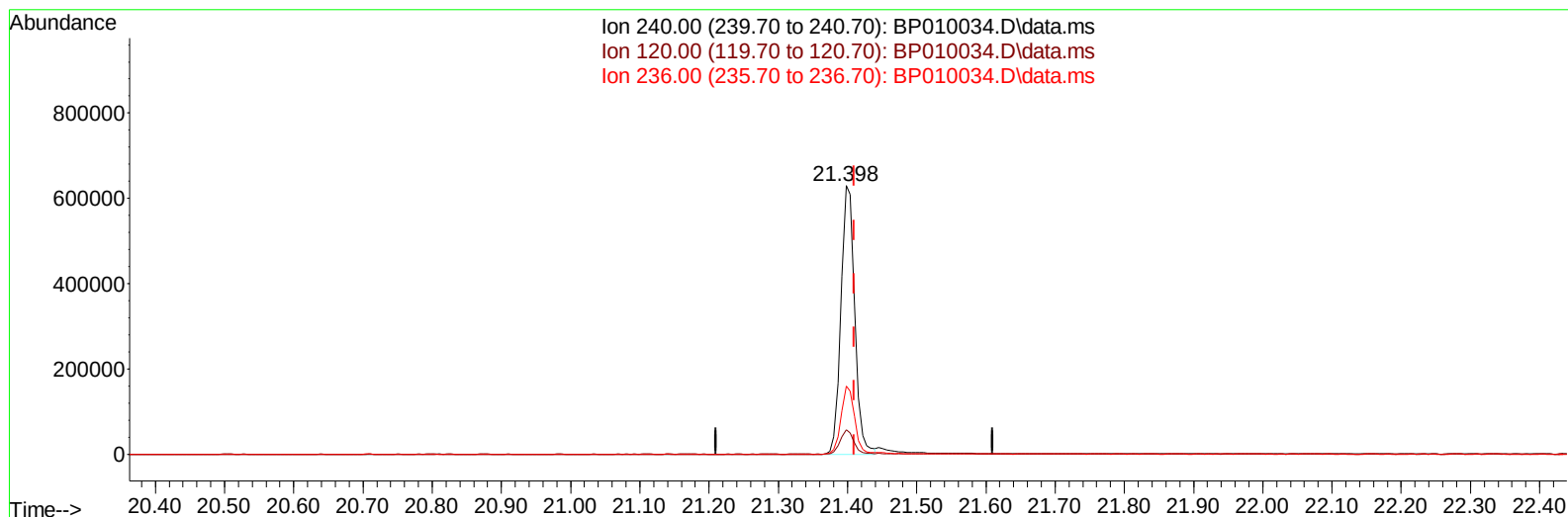
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010034.D
Acq On : 22 Apr 2022 02:18
Operator : CG/JU
Sample : N2473-16
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY9

Manual IntegrationsAPPROVED

Quant Time: Apr 22 04:23:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :mohammad ahmed 04/26/2022



TIC: BP010034.D\data.ms

(79) Chrysene-d12 (I)

21.398min (-0.012) 20.00 ng/u1

response 861002

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.14
236.00	31.60	25.28#
0.00	0.00	0.00

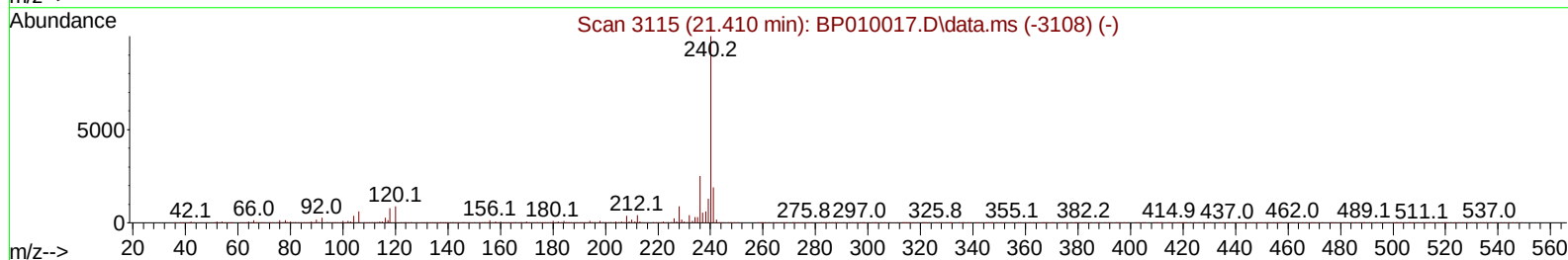
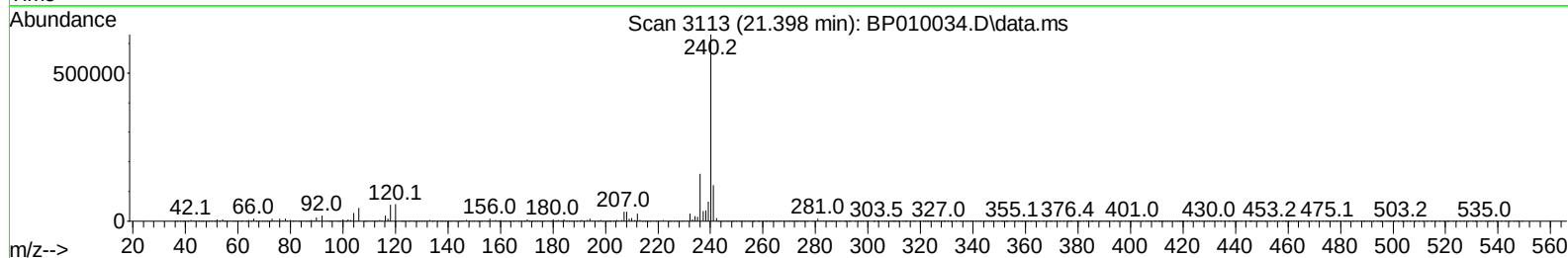
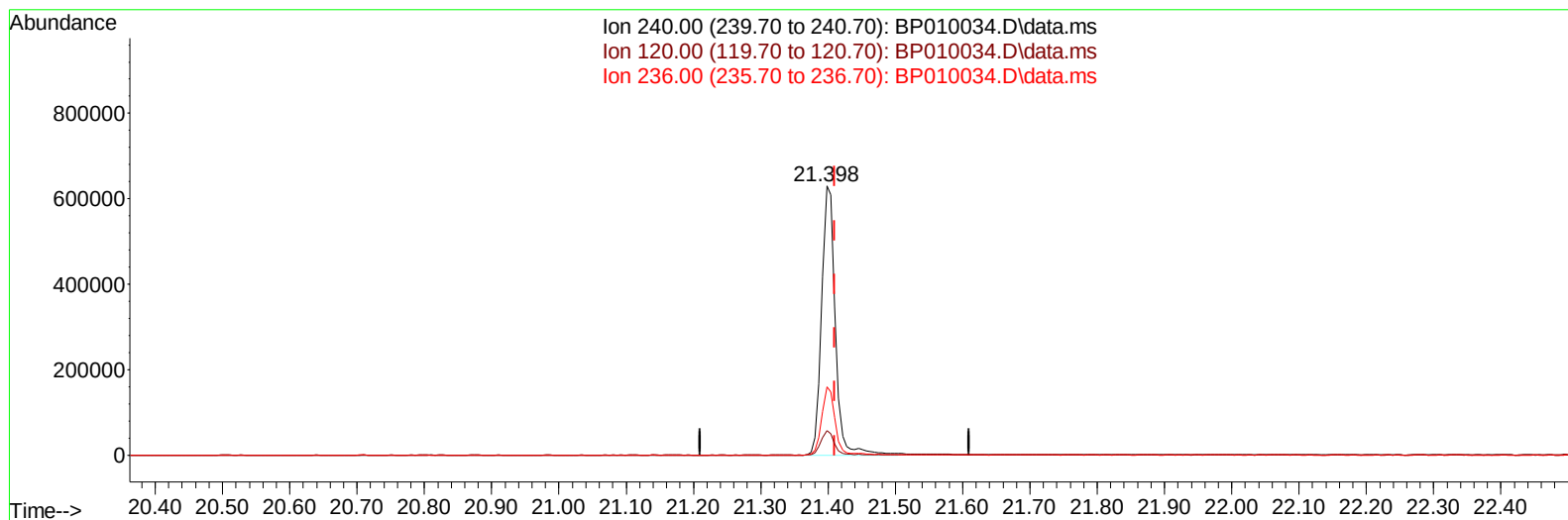
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(79) Chrysene-d12 (I)

21.398min (-0.012) 20.00 ng/ul m

response 889087

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.14
236.00	31.60	25.28#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	123733	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	572353	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	422511	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	946011	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	889087m	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264	800692	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	12981	4.575	ng/uL	0.00
4) Pyridine-d5	3.793	84	45025	5.641	ng/ul	0.00
7) Phenol-d5	7.099	99	65832	5.849	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	208548	30.881	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	190791	22.890	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	132785	14.240	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	131398	28.865	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	140094	28.179	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	237071	24.495	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	320060	23.063	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1046735	31.848	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1132176	28.463	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	29338	4.849	ng/ul	0.00
60) Fluorene-d10	15.563	176	928996	32.859	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	160912	26.900	ng/ul	0.00
73) Anthracene-d10	17.422	188	1621262	35.424	ng/ul	0.00
81) Pyrene-d10	19.651	212	1932826	38.226	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	1518590	36.177	ng/ul	0.00

Target Compounds Qvalue

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

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