

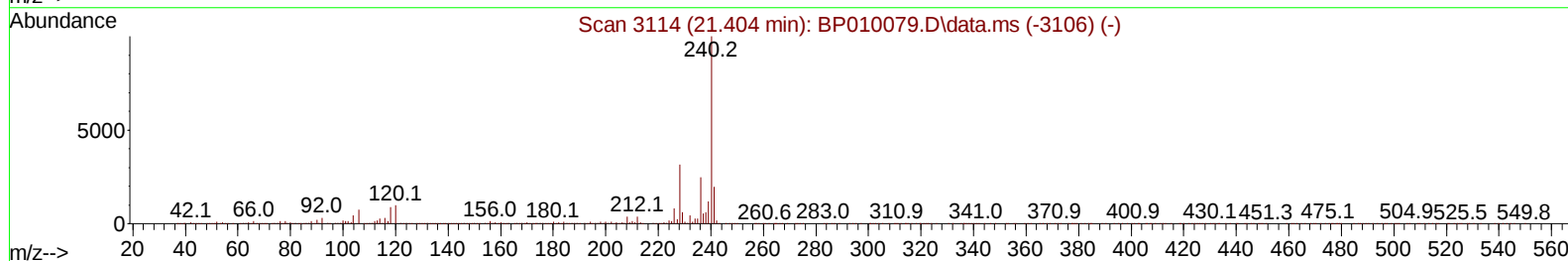
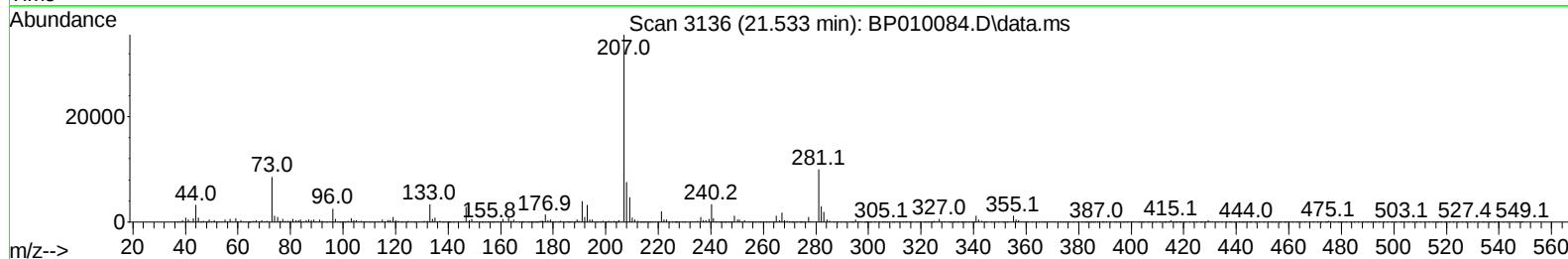
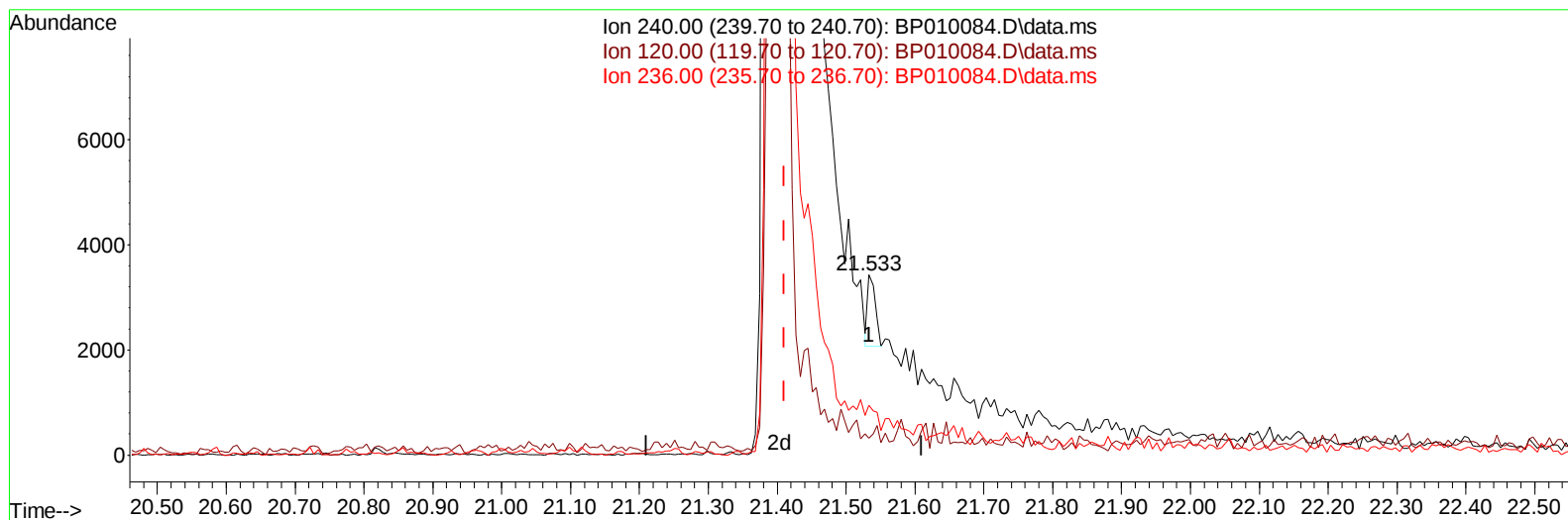
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010084.D
Acq On : 25 Apr 2022 15:58
Operator : CG/JU
Sample : N2475-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0BN2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/26/2022
Supervised By :Yogesh Patel 04/28/2022

Quant Time: Apr 26 01:02:28 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



TIC: BP010084.D\data.ms

(79) Chrysene-d12 (I)

21.533min (+ 0.123) 20.00 ng/u1

response 1074

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.39
236.00	31.60	27.51
0.00	0.00	0.00

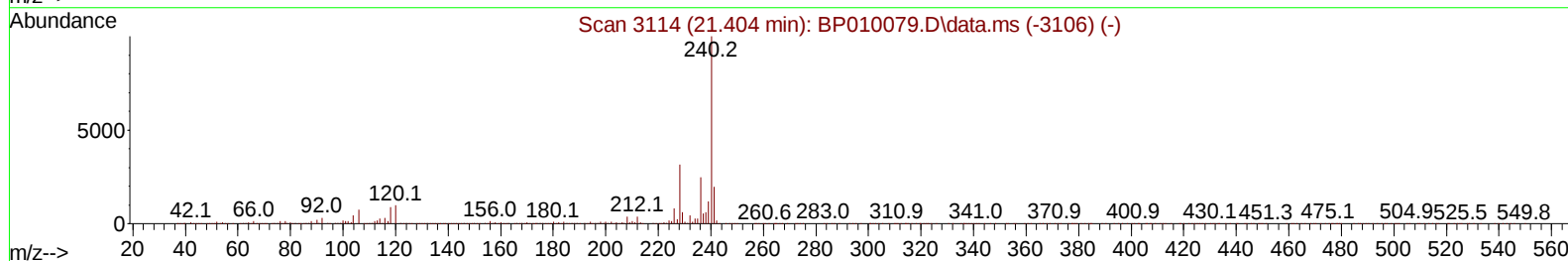
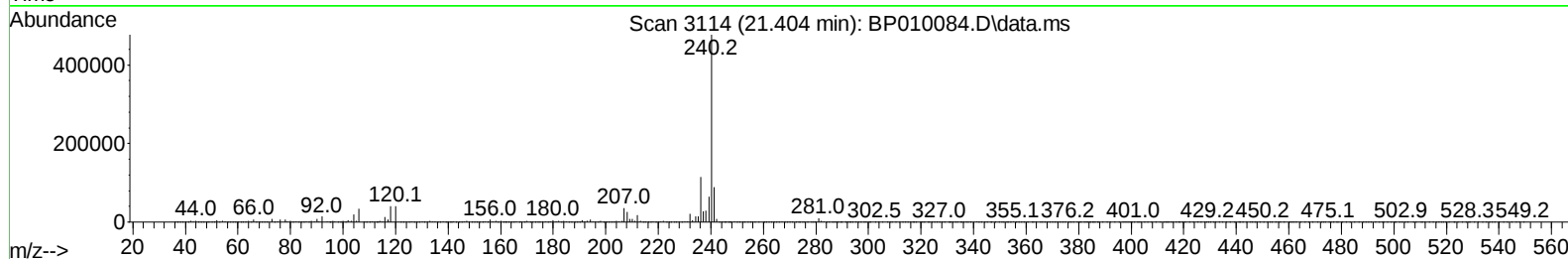
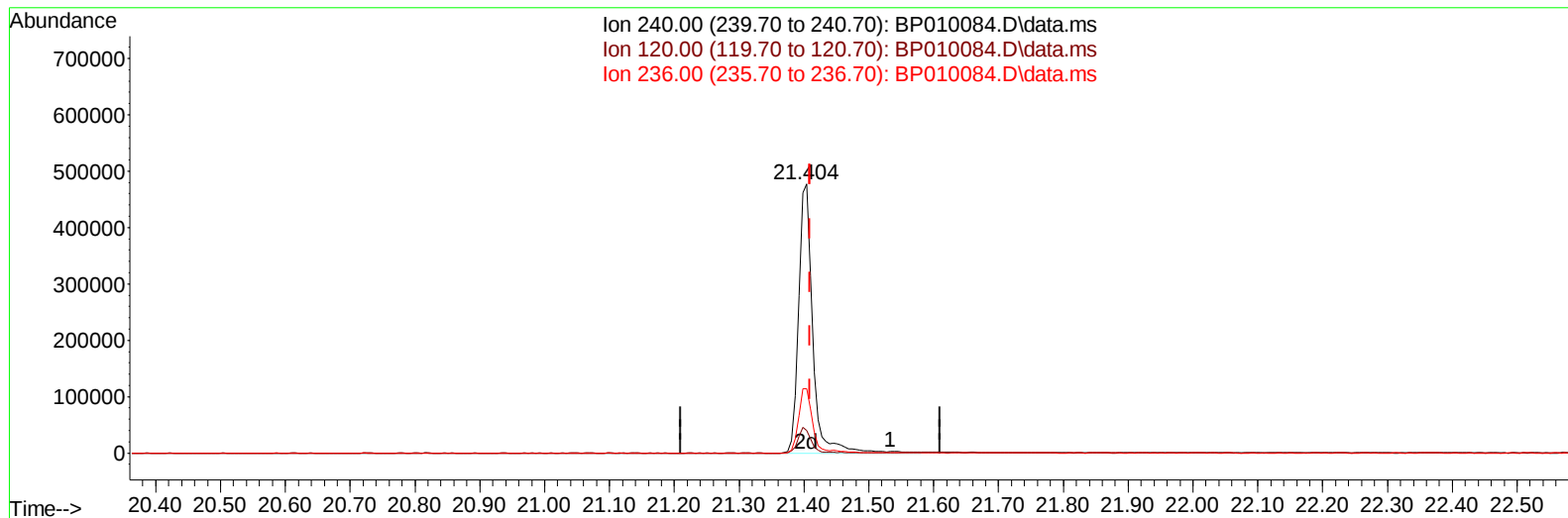
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TIC: BP010084.D\data.ms

(79) Chrysene-d12 (I)

21.404min (-0.006) 20.00 ng/ul m

response 726419

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.49
236.00	31.60	23.91#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	167160	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	711214	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	473509	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	954869	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	726419m	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	607461	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	11381	2.969	ng/uL	0.00
4) Pyridine-d5	3.799	84	32008	2.968	ng/ul	0.01
7) Phenol-d5	7.093	99	79657	5.239	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	263622	28.895	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	234556	20.830	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	156908	12.455	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	160092	28.302	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	163041	26.392	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	261380	21.733	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	329855	19.128	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1156866	31.408	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1235903	27.724	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	23676	3.492	ng/ul	0.00
60) Fluorene-d10	15.563	176	973271	30.718	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	160090	26.515	ng/ul	0.00
73) Anthracene-d10	17.416	188	1542599	33.393	ng/ul	0.00
81) Pyrene-d10	19.651	212	1649542	39.929	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	1084227	34.045	ng/ul	0.00

Target Compounds Qvalue

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

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