

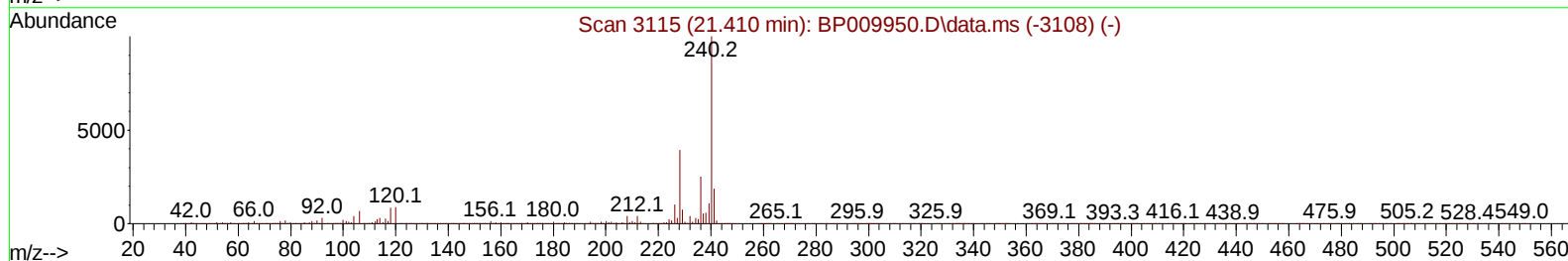
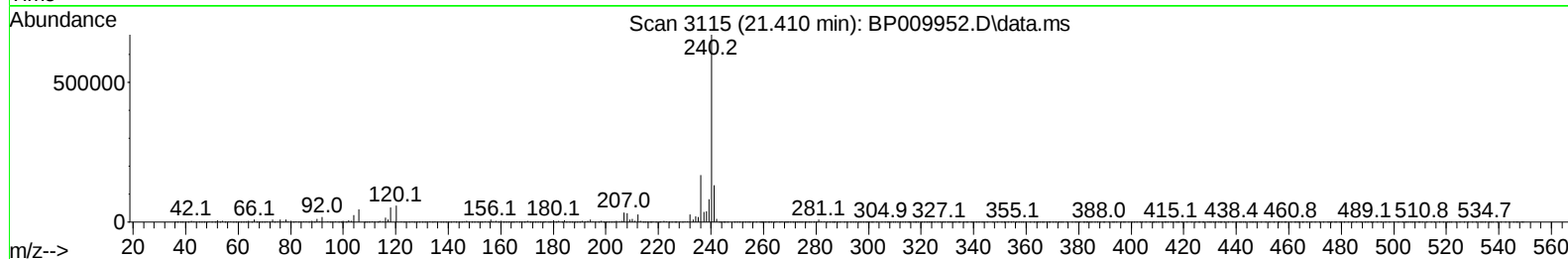
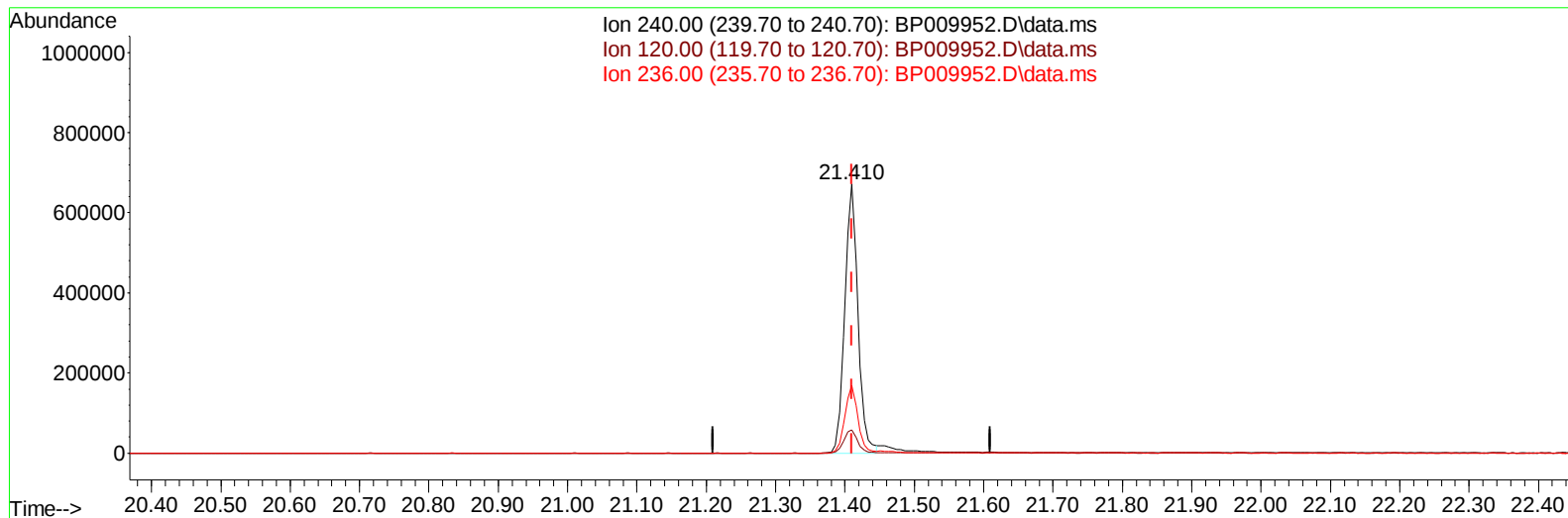
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009952.D
Acq On : 19 Apr 2022 19:10
Operator : CG/JU
Sample : PB144181BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK181

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/20/2022
Supervised By :Yogesh Patel 04/25/2022

Quant Time: Apr 20 01:53:16 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 20 01:32:14 2022
Response via : Initial Calibration



TIC: BP009952.D\data.ms

(79) Chrysene-d12 (I)

21.410min (+ 0.000) 20.00 ng/u1

response 874894

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.69
236.00	31.60	25.06#
0.00	0.00	0.00

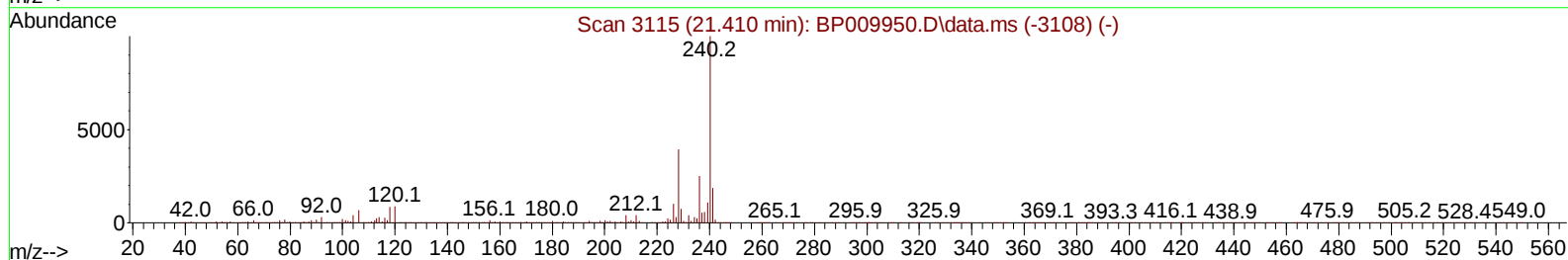
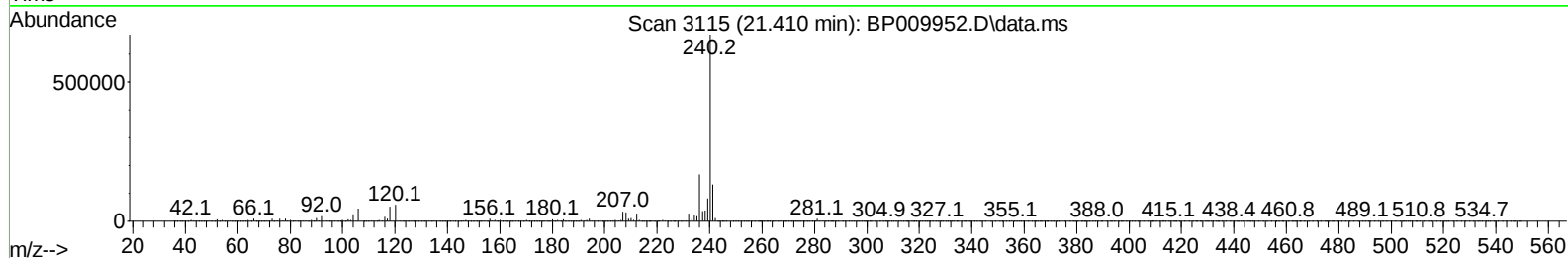
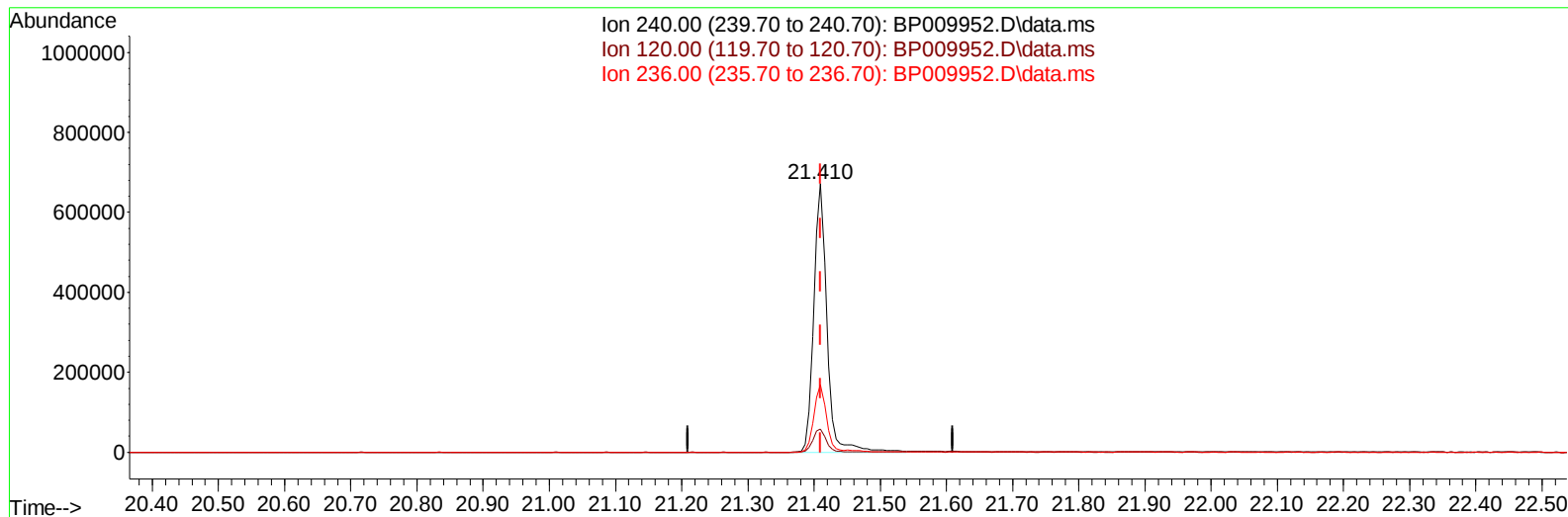
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(79) Chrysene-d12 (I)

21.410min (+ 0.000) 20.00 ng/ul m

response 920642

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.69
236.00	31.60	25.06#
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	7.946	152	142690	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	626962	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	445791	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	955240	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	920642m	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	825437	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	22416	7.011	ng/uL	0.00
4) Pyridine-d5	3.781	84	273253	30.998	ng/ul	0.00
7) Phenol-d5	7.099	99	379607	30.302	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	258707	34.689	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	310922	32.903	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	327817	31.347	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	174025	38.491	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	196622	39.667	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	346428	33.399	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	457841	30.991	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1215743	34.718	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1308977	31.552	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	179135	28.207	ng/ul	0.00
60) Fluorene-d10	15.569	176	1091849	37.065	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	175961	31.095	ng/ul	0.00
73) Anthracene-d10	17.428	188	1611767	35.241	ng/ul	0.00
81) Pyrene-d10	19.657	212	2115867	42.081	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1536060	35.734	ng/ul	0.00

Target Compounds Qvalue

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

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