

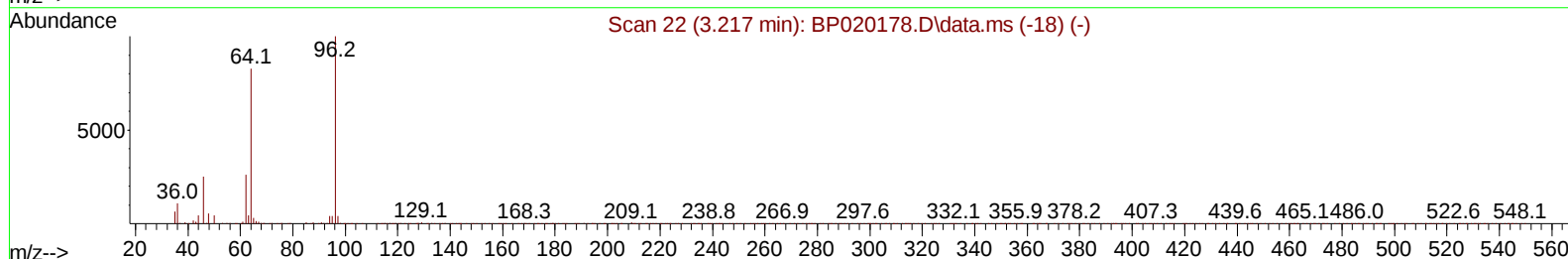
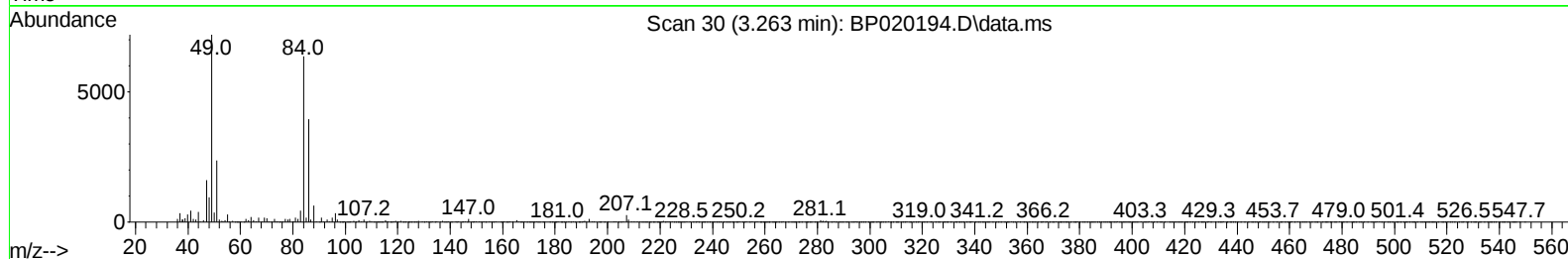
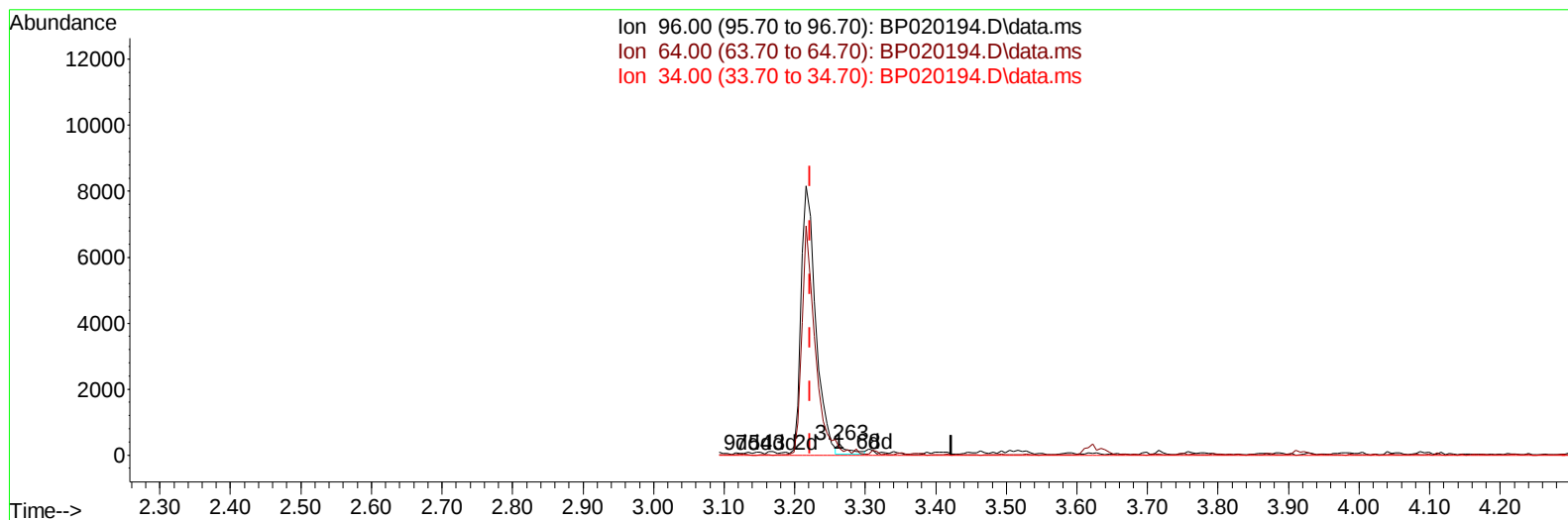
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
 Data File : BP020194.D
 Acq On : 04 May 2024 04:49
 Operator : MA/JU
 Sample : PB160586BL
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK586

Manual IntegrationsAPPROVED

Quant Time: May 04 05:36:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 03 06:07:25 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/06/2024
 Supervised By :mohammad ahmed 05/06/2024



TIC: BP020194.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.263min (+ 0.041) 0.14 ng/uL

response 321

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.80	61.31
34.00	0.00	0.00
0.00	0.00	0.00

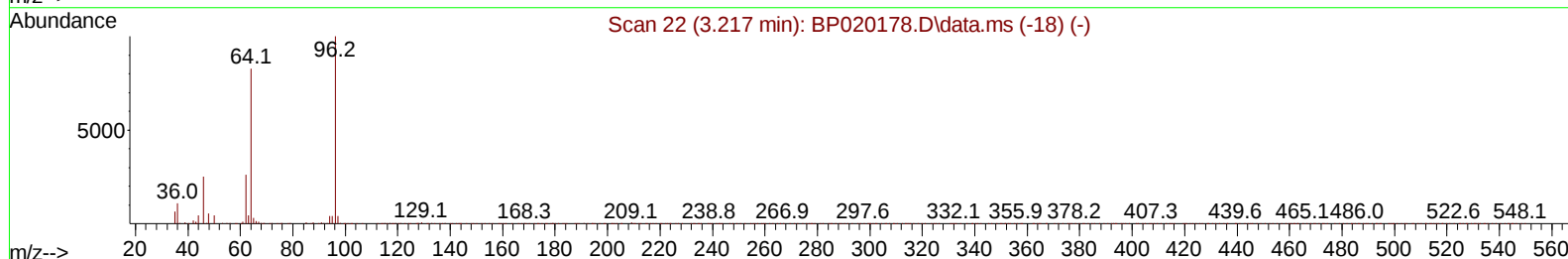
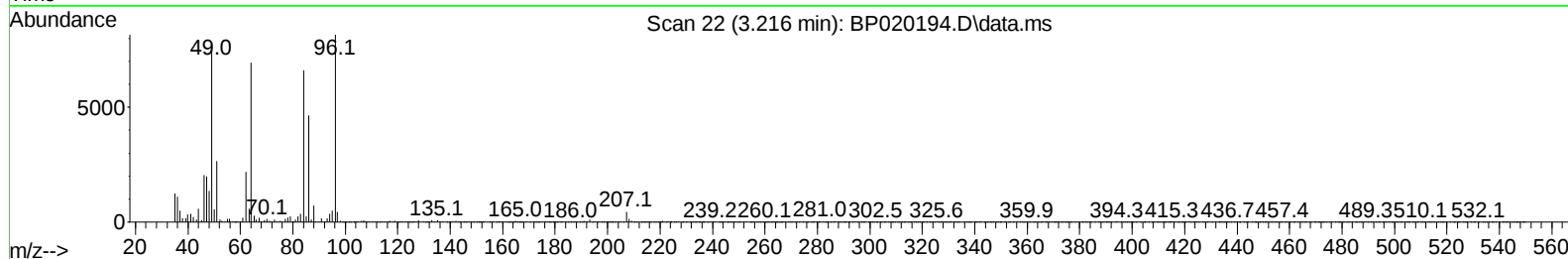
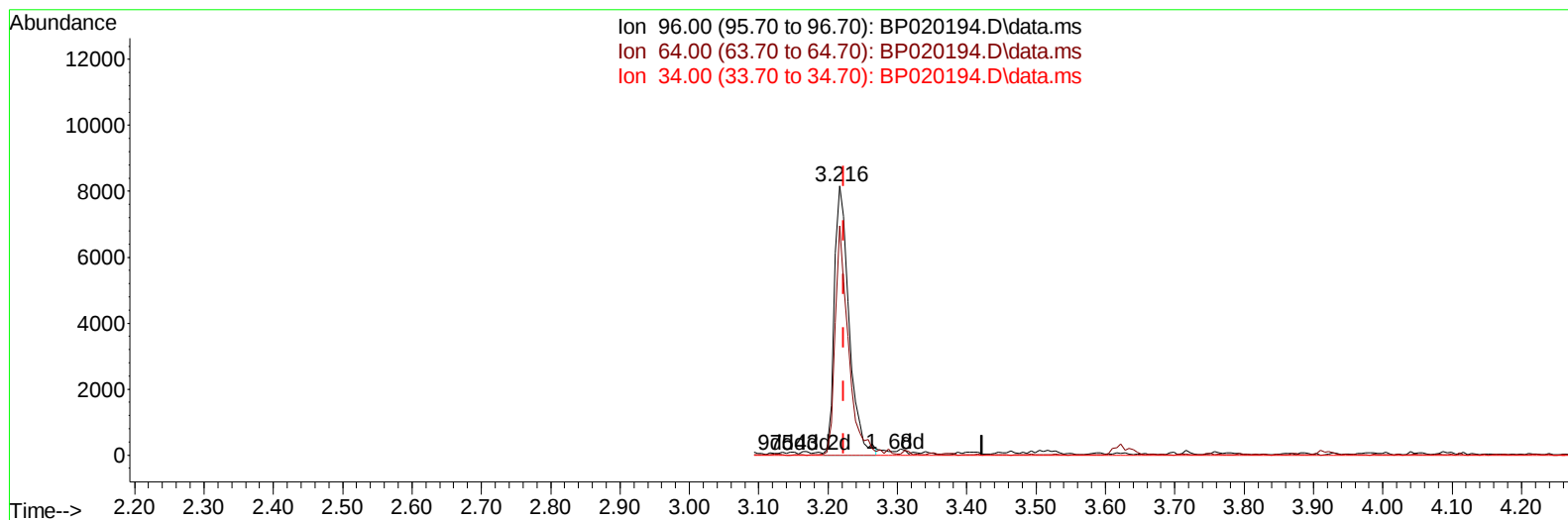
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TIC: BP020194.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.216min (-0.006) 5.39 ng/uL m

response 12032

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.80	85.25#
34.00	0.00	0.00
0.00	0.00	0.00

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Reviewed By :Yogesh Patel 05/06/2024
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Quant Time: May 04 05:36:37 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	94329	20.000	ng/ul	0.00
20) Naphthalene-d8	10.392	136	398124	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.286	164	304363	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.127	188	701862	20.000	ng/ul	0.00
79) Chrysene-d12	21.615	240	638503	20.000	ng/ul	0.00
88) Perylene-d12	24.962	264	592789	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	12032m	5.387	ng/uL	0.00
4) Pyridine-d5	3.616	84	160140	24.675	ng/ul	0.00
7) Phenol-d5	6.810	99	211419	26.098	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	136269	27.691	ng/ul	0.00
11) 2-Chlorophenol-d4	7.163	132	164149	28.345	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	175978	26.869	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	86155	28.899	ng/ul	0.00
24) 2-Nitrophenol-d4	9.492	143	101022	29.590	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	162363	26.061	ng/ul	0.00
31) 4-Chloroaniline-d4	10.551	131	251989	28.900	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	691293	31.102	ng/ul	-0.01
49) Acenaphthylene-d8	13.975	160	713925	29.077	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	86770	25.401	ng/ul	0.00
60) Fluorene-d10	15.316	176	565462	30.502	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	115694	25.976	ng/ul	0.00
73) Anthracene-d10	17.227	188	933510	30.745	ng/ul	-0.01
81) Pyrene-d10	19.627	212	1105838	29.628	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.733	264	846574	29.536	ng/ul	-0.02

Target Compounds Qvalue

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

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