

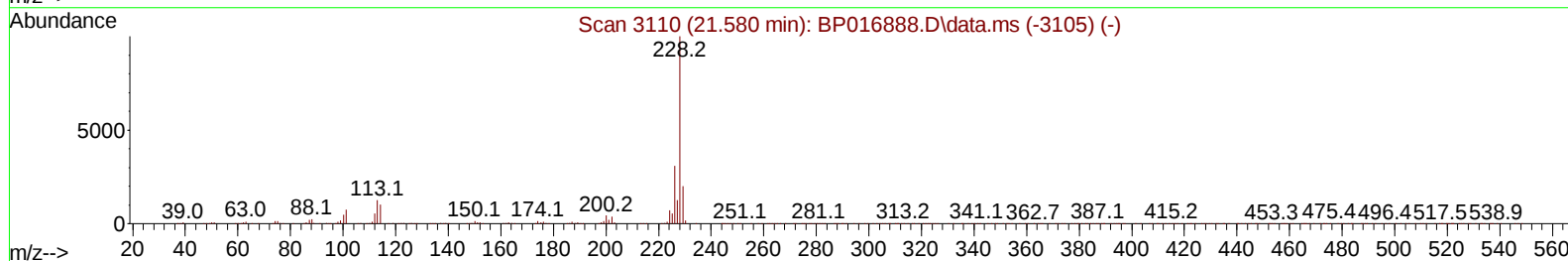
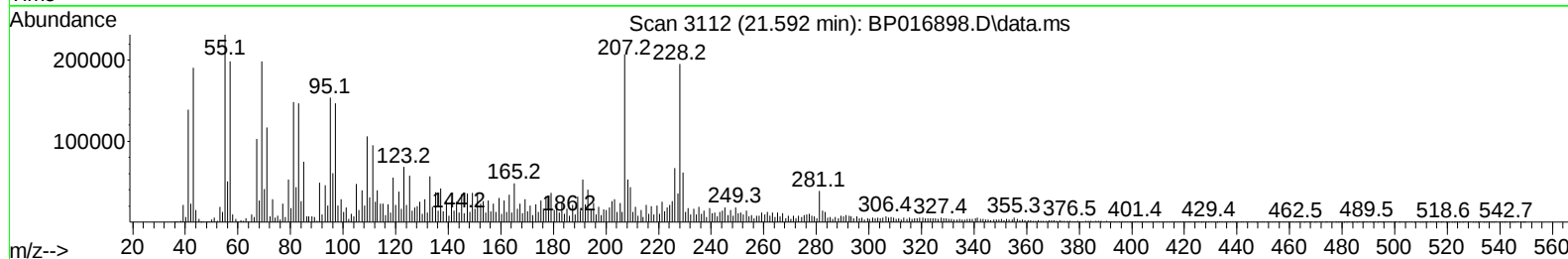
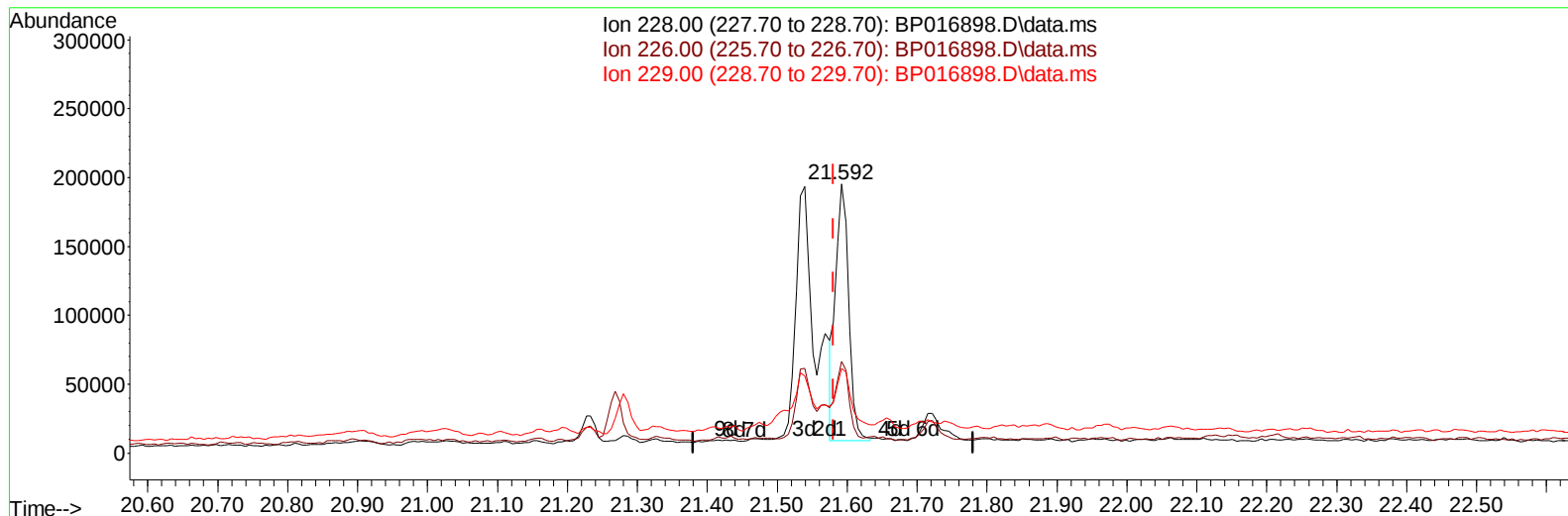
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083123\
Data File : BP016898.D
Acq On : 31 Aug 2023 15:18
Operator : MA/JU
Sample : 04113-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYH6

Manual IntegrationsAPPROVED

Quant Time: Sep 01 01:38:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 30 02:25:00 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/01/2023
Supervised By :mohammad ahmed 09/01/2023



TIC: BP016898.D\data.ms

(87) Chrysene

21.592min (+ 0.012) 2.77 ng/u1

response 245971

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	31.20	34.10
229.00	20.00	31.39#
0.00	0.00	0.00

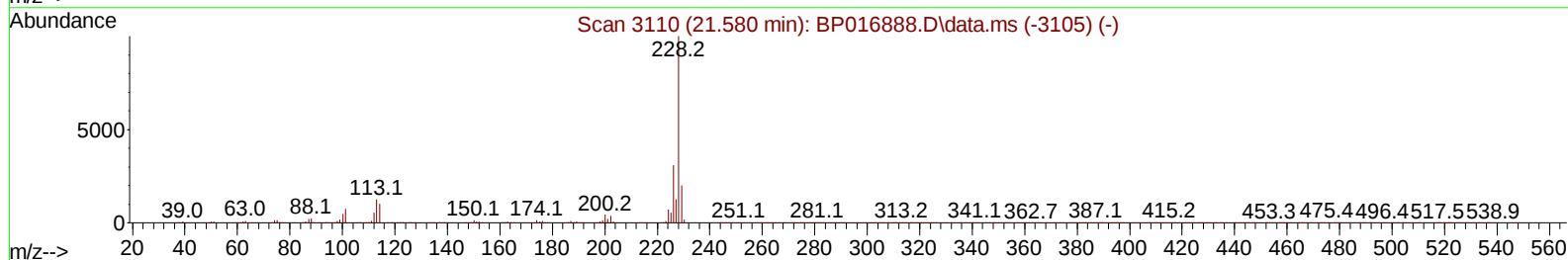
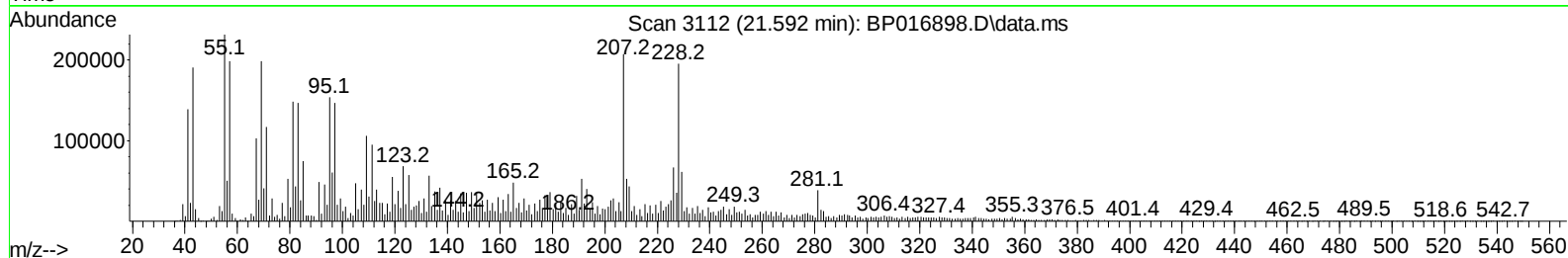
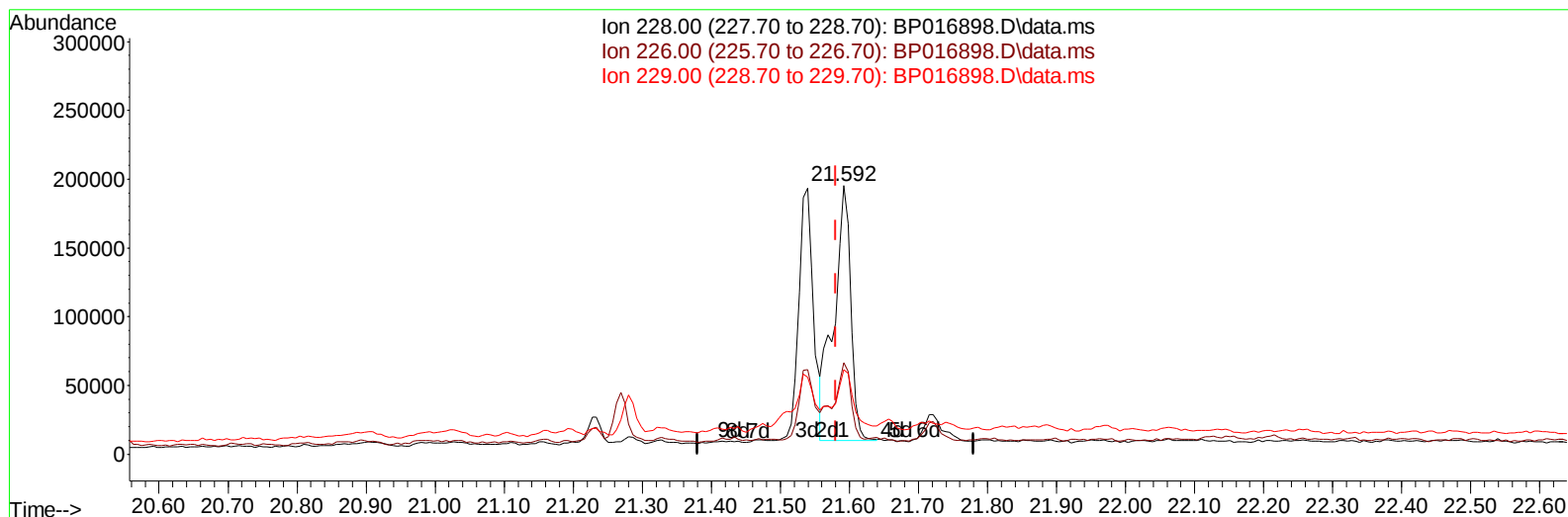
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(87) Chrysene

21.592min (+ 0.012) 3.59 ng/ul m

response 319591

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	31.20	34.10
229.00	20.00	31.39#
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	8.028	152	492950	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	1989431	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.681	164	1037169	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	1719546	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	1420970	20.000	ng/ul	0.02
88) Perylene-d12	24.133	264	1667383	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	36325	3.046	ng/uL	0.00
4) Pyridine-d5	3.828	84	33958	0.932	ng/ul	0.02
7) Phenol-d5	7.169	99	666417	15.203	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	449009	16.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	536219	16.270	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	357600	9.994	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	263657	17.528	ng/ul	0.00
24) 2-Nitrophenol-d4	9.940	143	266143	16.786	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	446761	15.114	ng/ul	0.00
31) 4-Chloroaniline-d4	11.016	131	144319	3.184	ng/ul	0.00
46) Dimethylphthalate-d6	14.092	166	1369213	17.505	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	1549135	17.520	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	123802	7.658	ng/ul	0.01
60) Fluorene-d10	15.675	176	1106160	16.794	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	18069	1.764	ng/ul	0.01
73) Anthracene-d10	17.551	188	1410319	18.503	ng/ul	0.00
81) Pyrene-d10	19.786	212	1436285	18.404	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.962	264	1507010	18.043	ng/ul	0.02
Target Compounds						
					Qvalue	
16) Acetophenone	8.875	105	165507	2.983	ng/ul	99
30) Naphthalene	10.904	128	205719	1.996	ng/ul	99
36) 2-Methylnaphthalene	12.510	142	115429	1.623	ng/ul	100
72) Phenanthrene	17.492	178	317890	3.512	ng/ul	99
80) Fluoranthene	19.457	202	485858	5.183	ng/ul	98
82) Pyrene	19.816	202	554689	5.728	ng/ul	100
85) Benzo(a)anthracene	21.539	228	270296	2.762	ng/ul#	89
87) Chrysene	21.592	228	319591m	3.595	ng/ul	
90) Benzo(b)fluoranthene	23.327	252	478738	4.496	ng/ul#	72
91) Benzo(k)fluoranthene	23.380	252	126233	1.199	ng/ul#	47
93) Benzo(a)pyrene	24.015	252	301091	3.126	ng/ul#	70
94) Indeno(1,2,3-cd)pyrene	26.868	276	230257	2.340	ng/ul	92
96) Benzo(g,h,i)perylene	27.727	276	246977	3.310	ng/ul#	88

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

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