

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021920\

Data File : BP001789.D

Acq On : 19 Feb 2020 21:43

Operator : CG/JU

Sample : L1424-11

Misc :

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Feb 20 01:29:43 2020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Feb 19 02:37:06 2020

Response via : Initial Calibration

Instrument :

BNA_P

ClientSampleId :

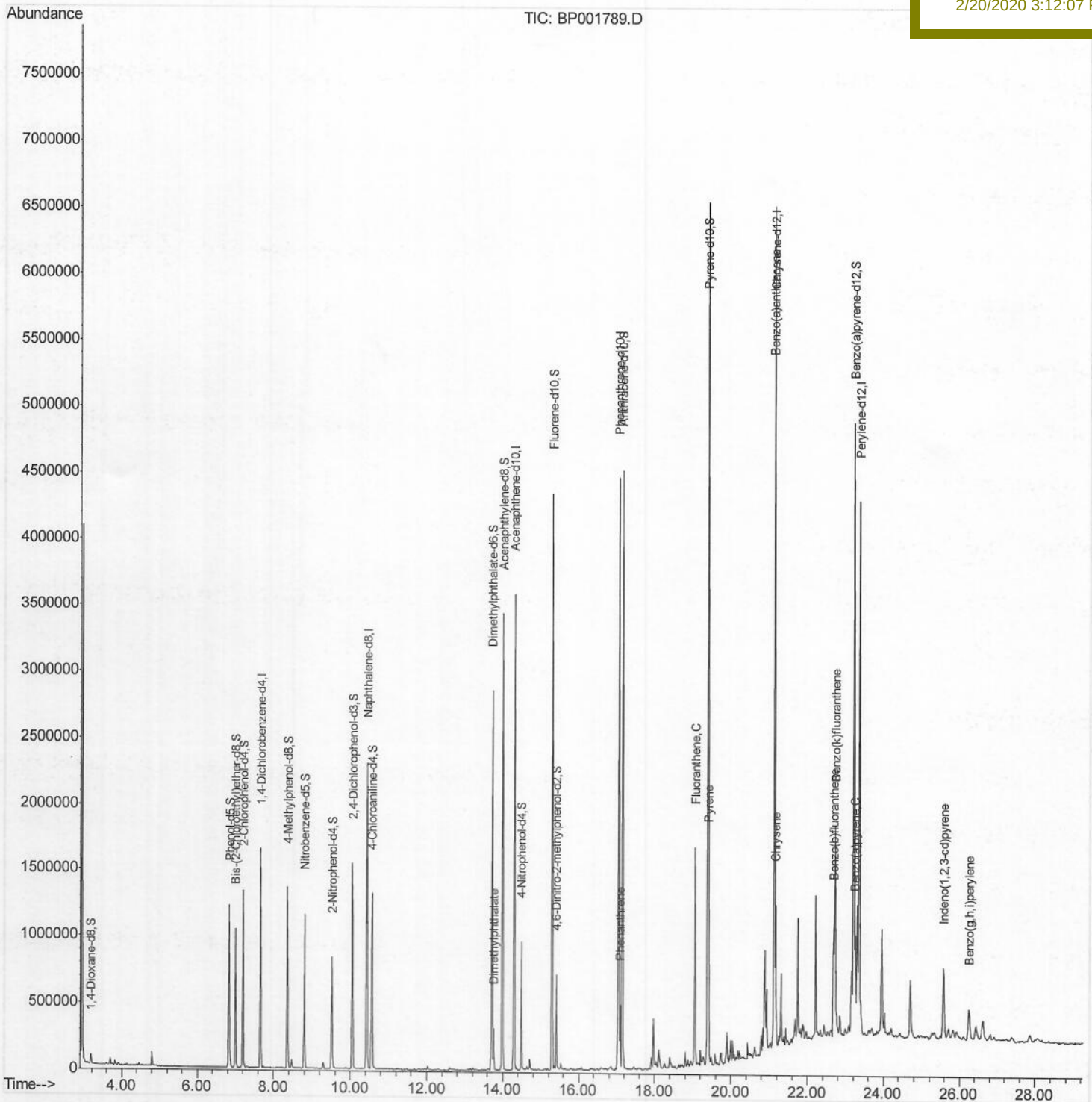
C5B11

Manual Integrations

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mohammad

2/20/2020 3:12:07 PM



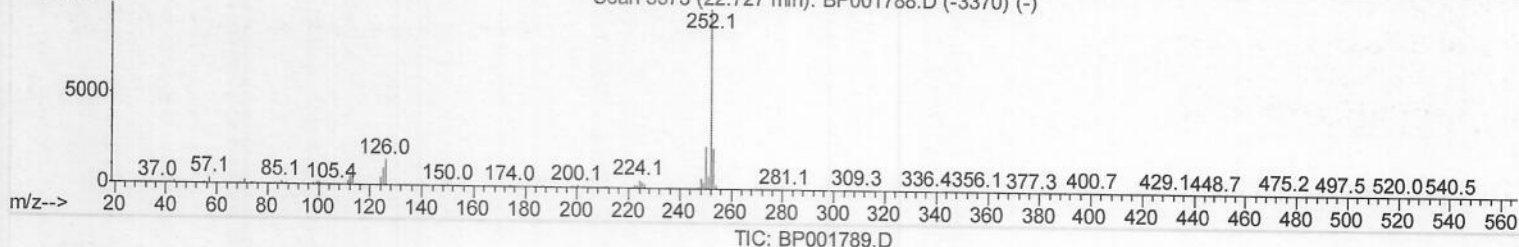
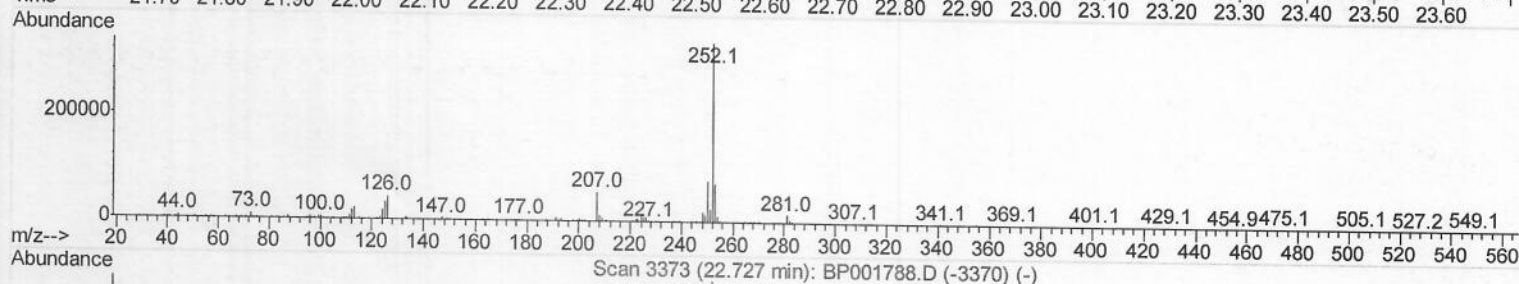
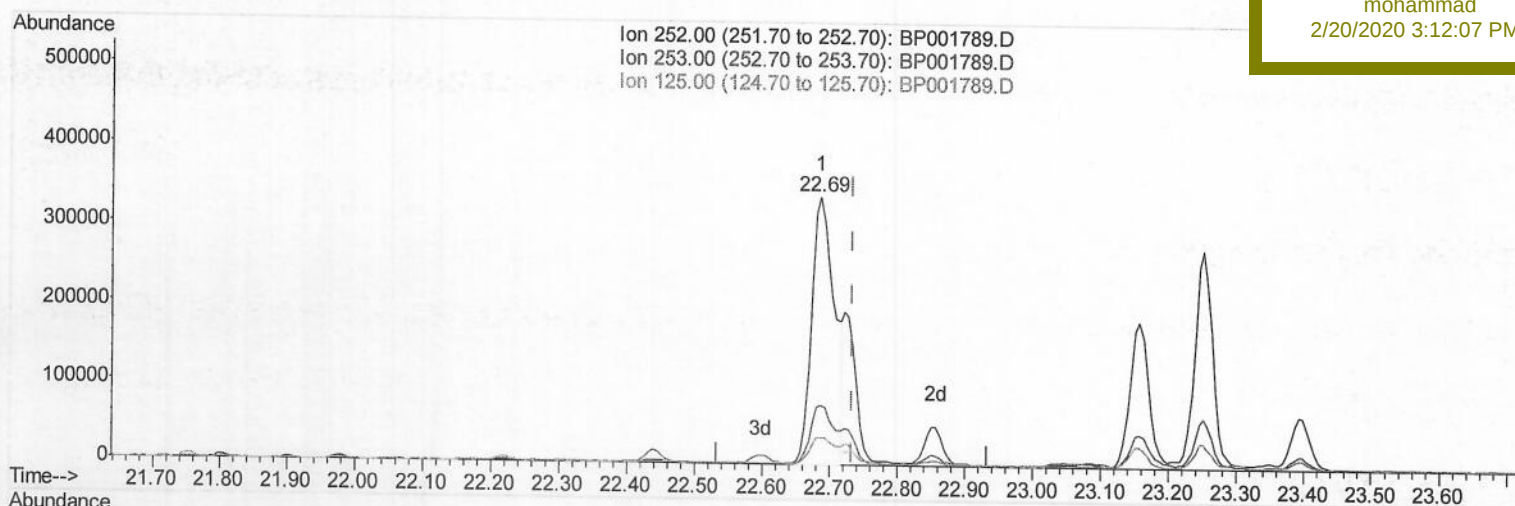
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(89) Benzo(k)fluoranthene

22.686min (-0.047) 4.37ng/ul

response 783870

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	22.17
125.00	10.00	10.12
0.00	0.00	0.00

Quantitation Report (Qedit)

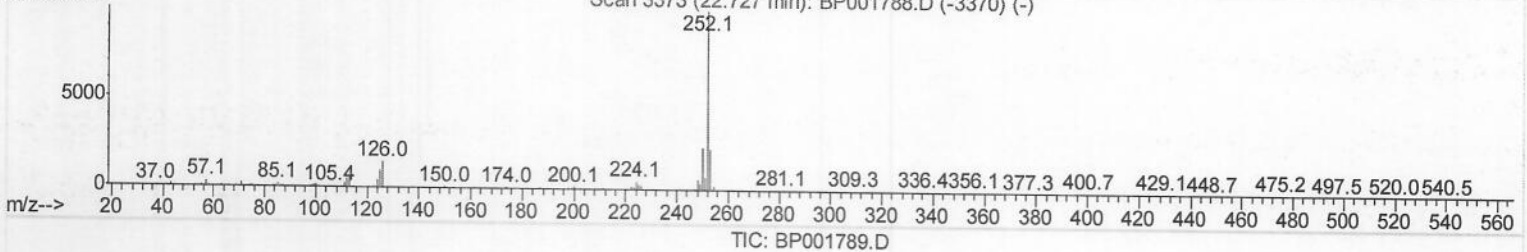
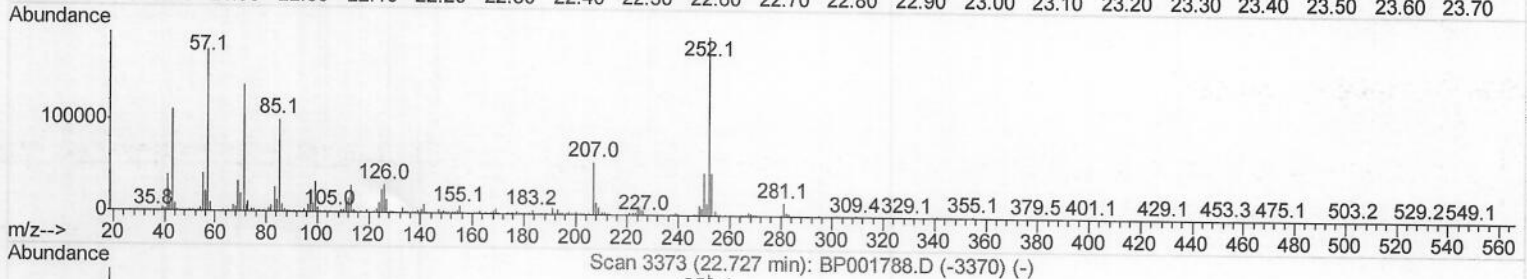
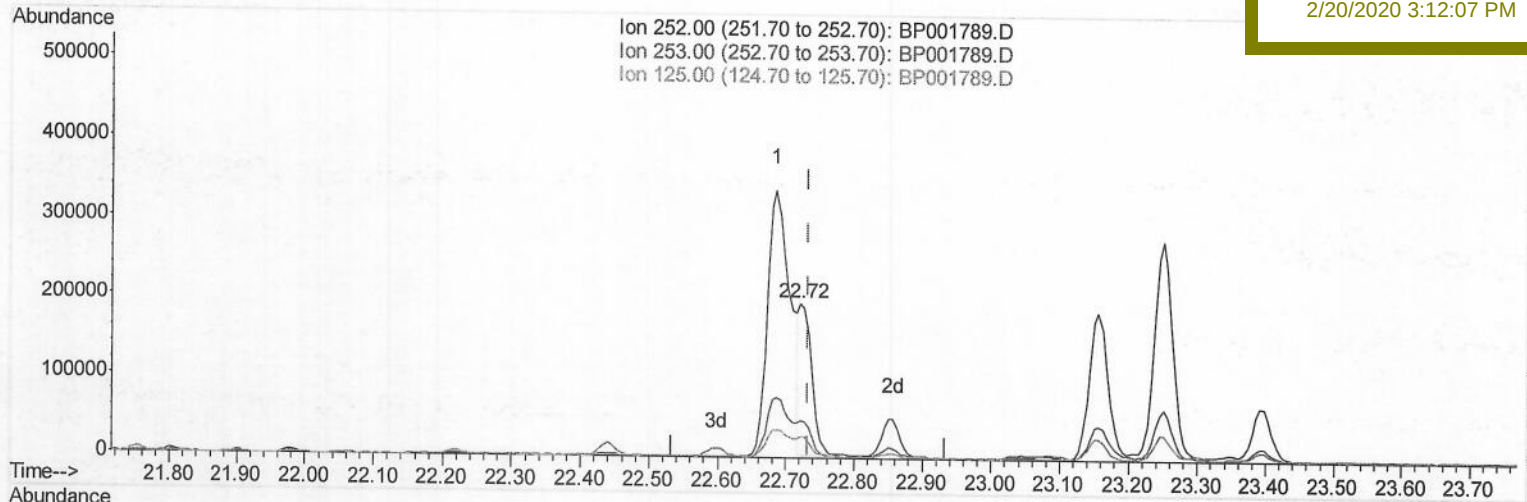
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 QLast Update : Wed Feb 19 02:37:06 2020
 Response via : Initial Calibration

Instrument :
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 C5B11

Manual Integrations
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 2/20/2020 3:12:07 PM



(89) Benzo(k)fluoranthene

22.721min (-0.012) 1.37ng/ul m > JV 02/25/20

response 245594

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	23.83
125.00	10.00	13.31#
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	465538	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1983005	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	1235542	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	2669175	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	2759119	20.00	ng/ul	0.00
86) Perylene-d12	23.35	264	2940019	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	37321	3.33	ng/uL	0.00
5) Phenol-d5	6.83	99	709088	20.16	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	437462	21.41	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	618431	21.75	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	516911	18.53	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	289292	21.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	289695	23.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	609109	21.10	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	734163	21.50	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	1933665	22.30	ng/ul	0.00
47) Acenaphthylene-d8	13.98	160	2367667	22.18	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	239918	15.89	ng/ul	0.00
58) Fluorene-d10	15.29	176	1773506	22.74	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	162910	15.48	ng/ul	0.00
71) Anthracene-d10	17.14	188	2599038	22.10	ng/ul	0.00
79) Pyrene-d10	19.39	212	3091707	21.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.21	264	3330991	23.46	ng/ul	0.00

Target Compounds

					Ovalue
45) Dimethylphthalate	13.75	163	195624	2.144	ng/ul 100
70) Phenanthrene	17.09	178	295872	2.010	ng/ul 99
77) Fluoranthene	19.06	202	943152	5.983	ng/ul 99
80) Pyrene	19.42	202	856545	4.564	ng/ul 98
83) Benzo(a)anthracene	21.12	228	564763	3.108	ng/ul 100
85) Chrysene	21.17	228	569194	3.207	ng/ul 97
88) Benzo(b)fluoranthene	22.69	252	783870	4.531	ng/ul 100
89) Benzo(k)fluoranthene	22.72	252	245594m	1.370	ng/ul 99
91) Benzo(a)pyrene	23.25	252	512262	3.160	ng/ul 98
92) Indeno(1,2,3-cd)pyrene	25.57	276	335453	1.656	ng/ul 96
94) Benzo(a,h,i)perylene	26.26	276	284645	1.653	ng/ul 99

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

JO 02/25/20