

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Quantitation Report (QT Reviewed)

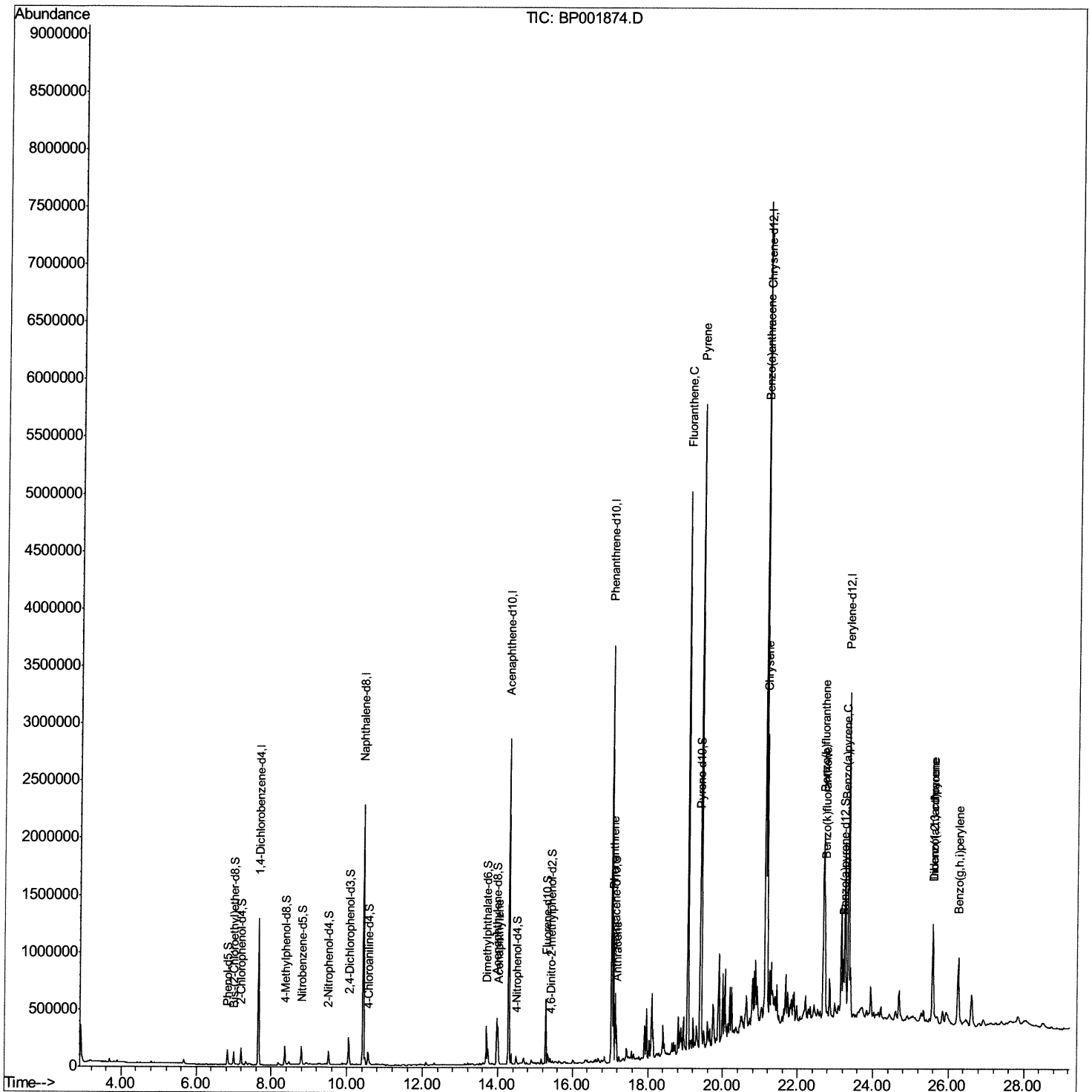
Data File : BP001874.D
Acq On : 24 Feb 2020 22:49
Operator : CG/JU
Sample : L1418-20DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C5AW4DL

Manual Integrations
APPROVED

mohammad
2/25/2020 3:42:35 PM

Quant Time: Feb 25 06:38:33 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 25 05:55:24 2020
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Quantitation Report (Qedit)

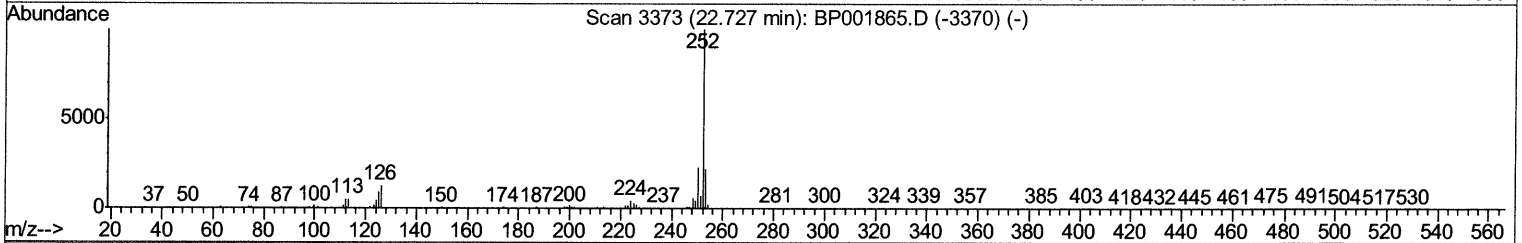
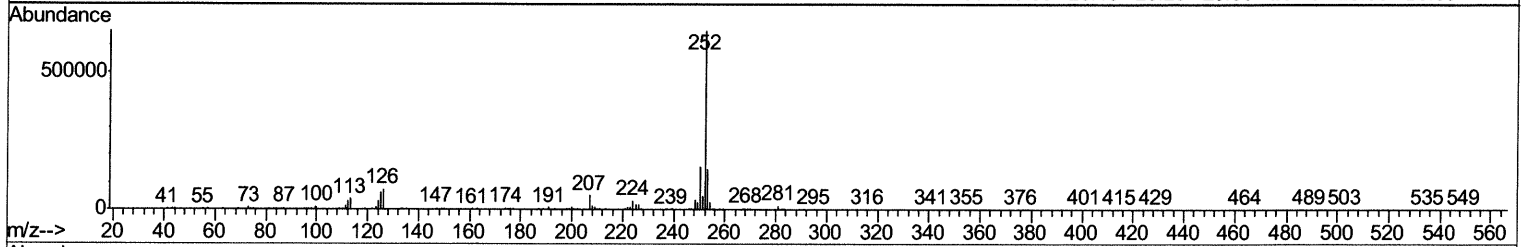
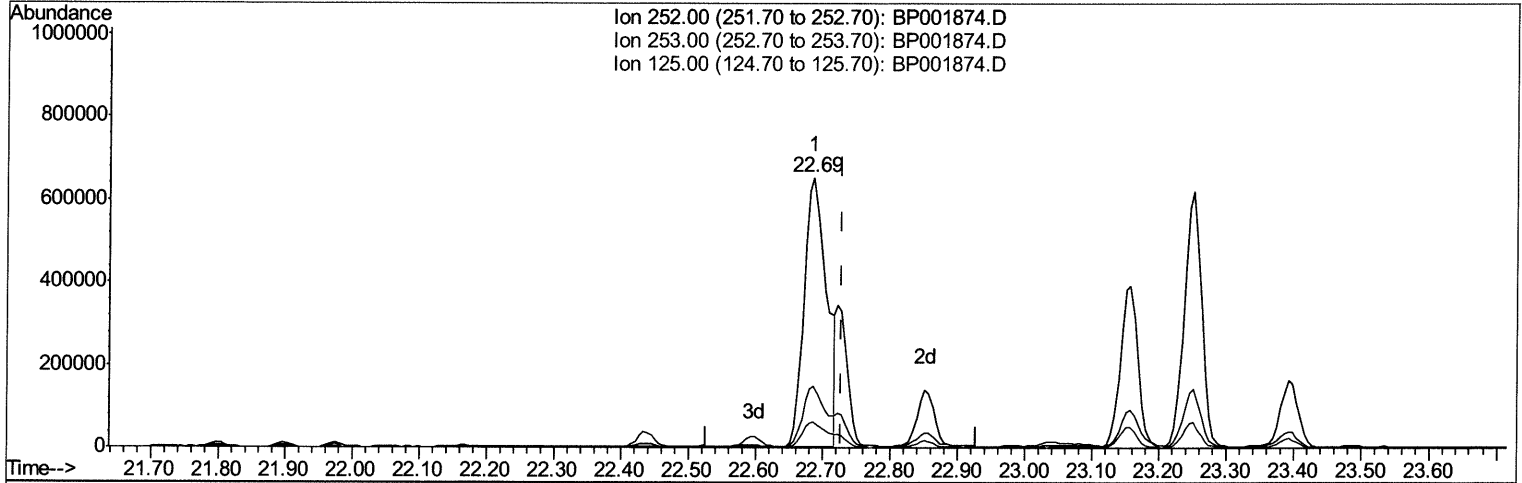
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TIC: BP001874.D

(89) Benzo(k)fluoranthene
22.686min (-0.041) 11.92ng/ul
response 1513759

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	22.78
125.00	10.00	9.40
0.00	0.00	0.00

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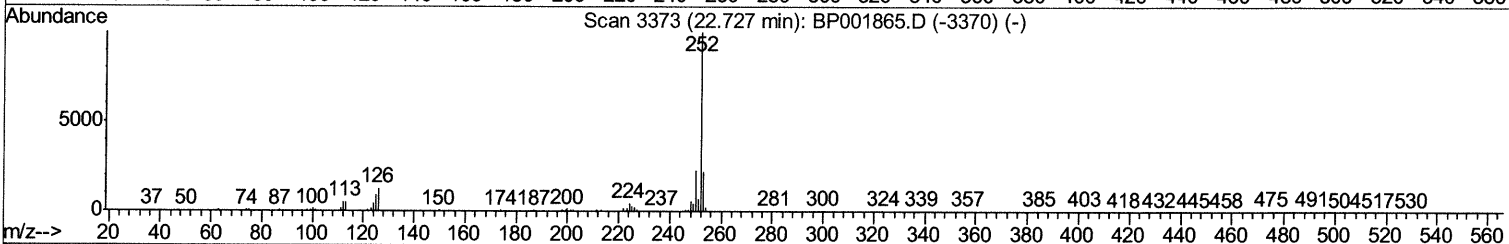
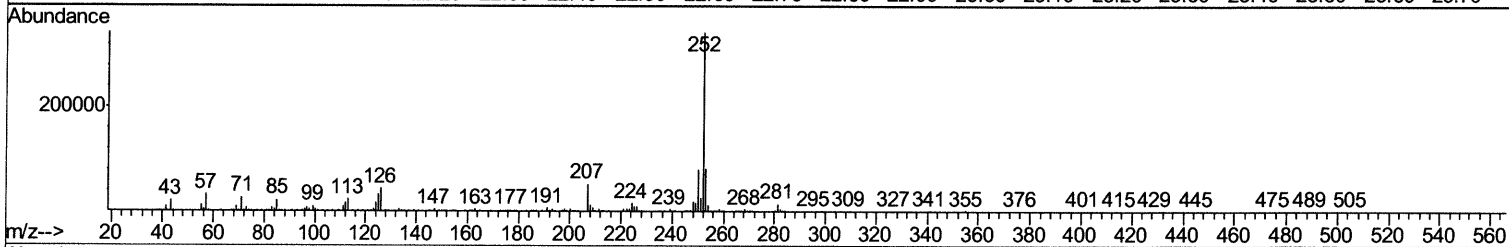
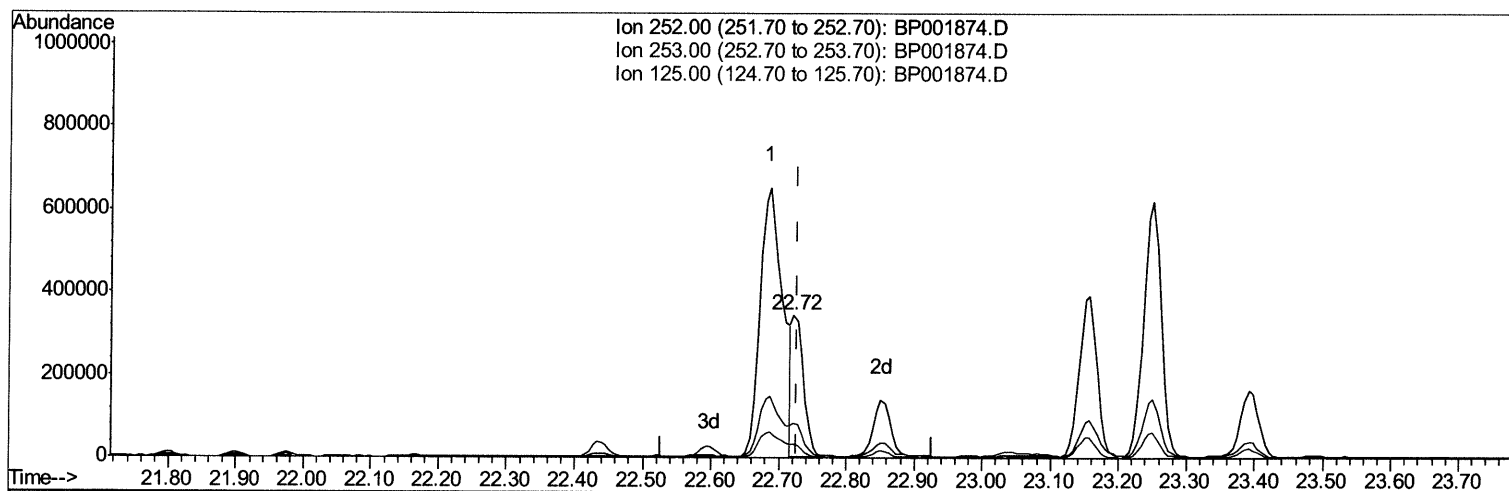
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TIC: BP001874.D

(89) Benzo(k)fluoranthene

22.721min (-0.006) 3.18ng/ul m *Juoz/27/20*

response 404216

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	24.22
125.00	10.00	9.36
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	369247	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	1968466	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	993916	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	2213999	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	2832107	20.00	ng/ul	0.00
86) Perylene-d12	23.34	264	2082819	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	5261	0.59	ng/uL	0.00
5) Phenol-d5	6.83	99	79010	2.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	46484	2.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	72198	3.20	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	64406	2.91	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	41847	3.14	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	42547	3.53	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	96979	3.38	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	73230	2.16	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	240506	3.45	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	285309	3.32	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	22733	1.87	ng/ul	0.00
58) Fluorene-d10	15.29	176	220733	3.52	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	13915	1.59	ng/ul	0.00
71) Anthracene-d10	17.14	188	352616	3.62	ng/ul	0.00
79) Pyrene-d10	19.38	212	555788	3.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.20	264	368900	3.67	ng/ul	0.00

Target Compounds

						Qvalue
48) Acenaphthylene	14.00	152	245772	2.638	ng/ul	100
70) Phenanthrene	17.08	178	679979	5.569	ng/ul	100
72) Anthracene	17.17	178	203438	1.687	ng/ul	99
77) Fluoranthene	19.06	202	2963566	22.666	ng/ul	99
80) Pyrene	19.41	202	3418919	17.748	ng/ul	100
83) Benzo(a)anthracene	21.12	228	1529818	8.203	ng/ul	97
85) Chrysene	21.17	228	1403416	7.703	ng/ul	98
88) Benzo(b)fluoranthene	22.69	252	1513759	12.351	ng/ul	99
89) Benzo(k)fluoranthene	22.72	252	404216m >	3.182	ng/ul >	94
91) Benzo(a)pyrene	23.25	252	1116972	9.725	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.57	276	730737	5.093	ng/ul#	94
93) Dibenzo(a,h)anthracene	25.59	278	152161	1.263	ng/ul#	81
94) Benzo(g,h,i)perylene	26.26	276	727361	5.964	ng/ul	97

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

Jul 27/20