

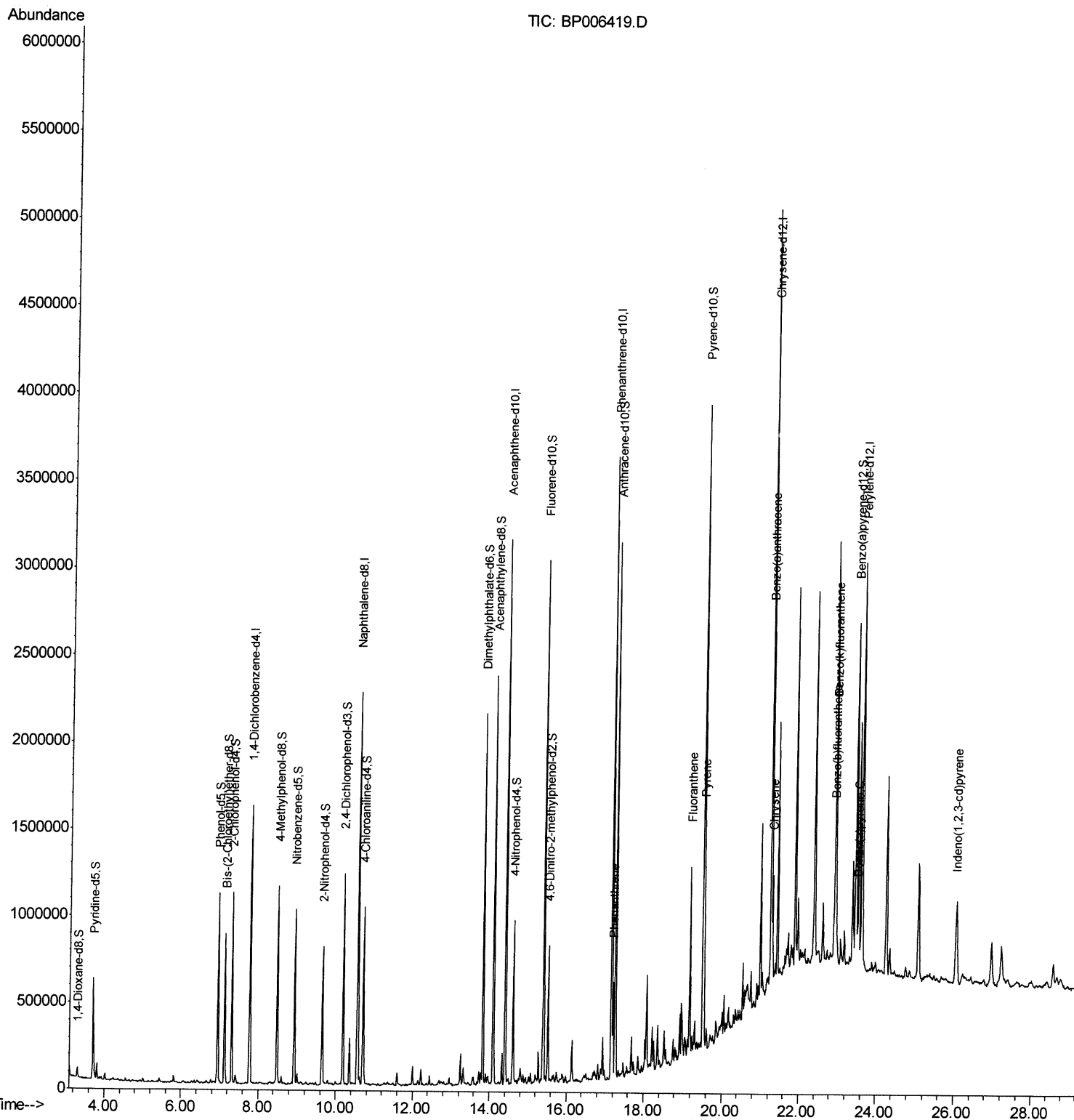
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP072921\
Data File : BP006419.D
Acq On : 29 Jul 2021 14:17
Operator : CG/JU
Sample : M3123-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0080

Quant Time: Jul 29 15:04:04 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 29 14:34:44 2021
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
7/30/2021 4:30:33 PM



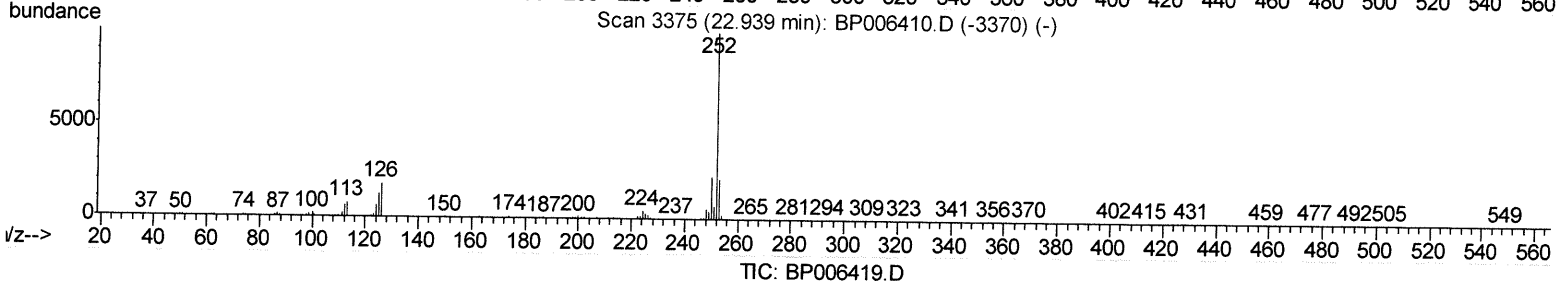
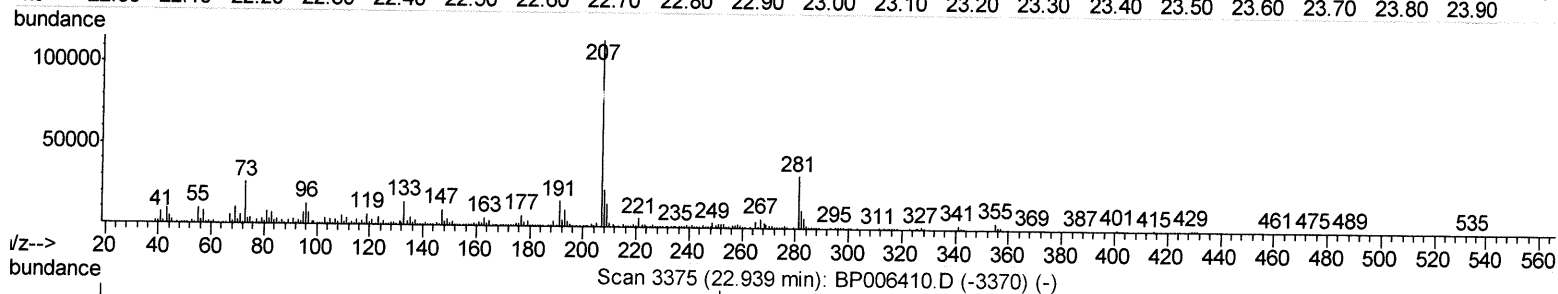
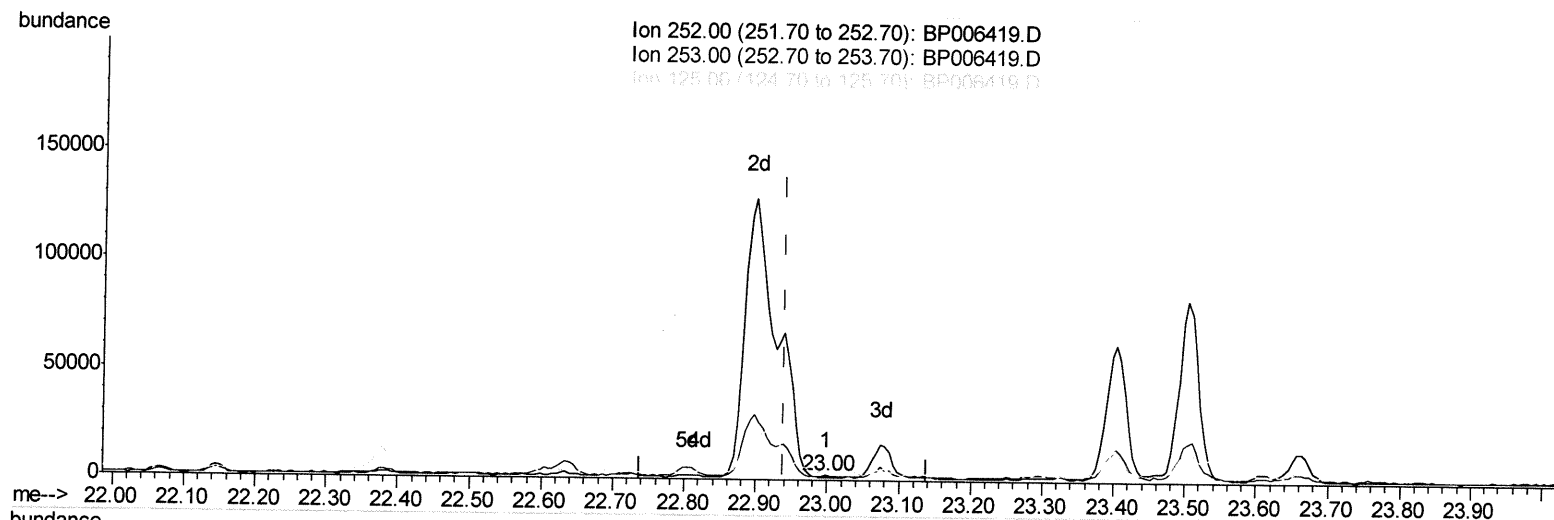
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(91) Benzo(k)fluoranthene

22.998min (+0.059) 0.01ng/ul

response 658

Ion	Exp%	Act%
252.00	100	100
253.00	21.80	94.19#
125.00	12.80	126.44#
0.00	0.00	0.00

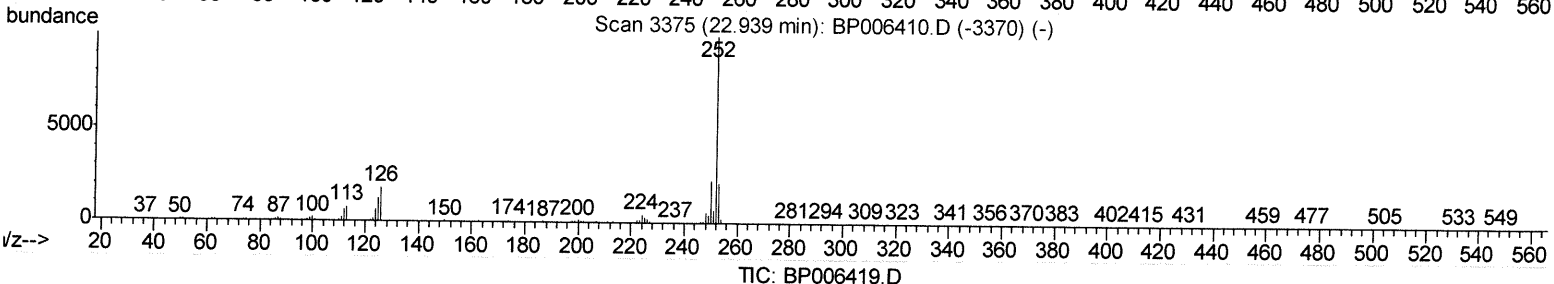
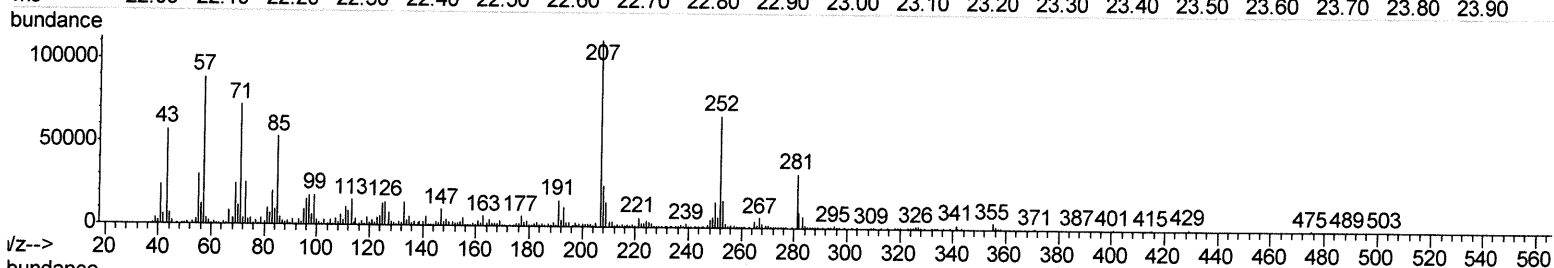
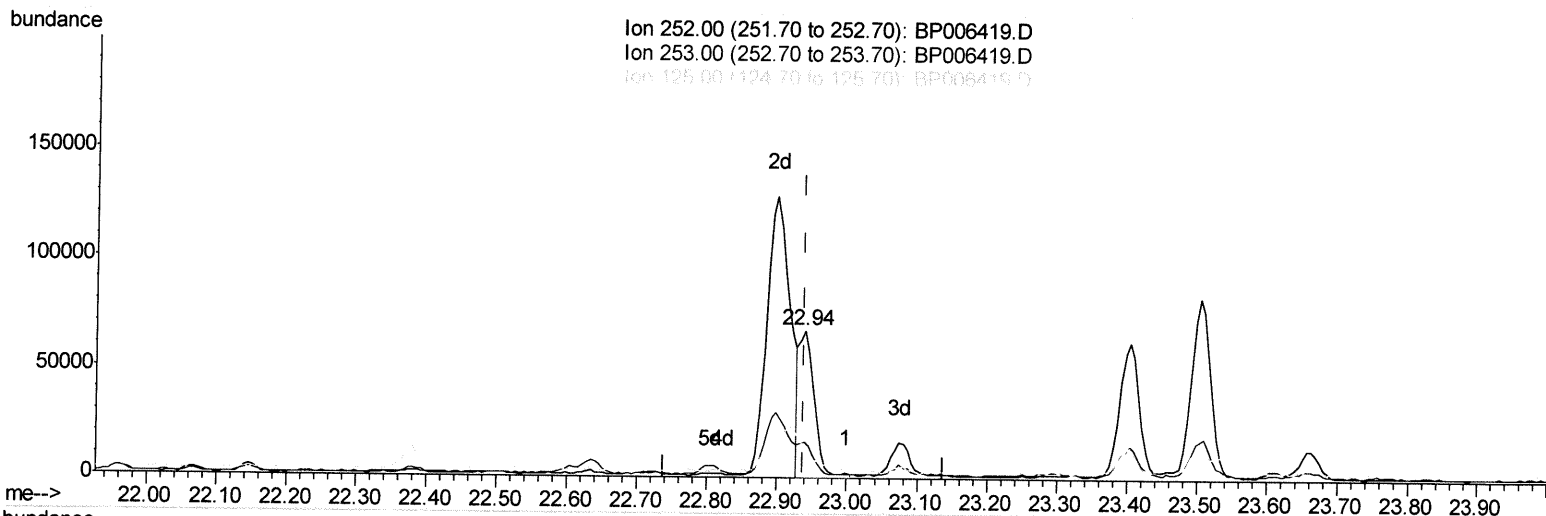
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(91) Benzo(k)fluoranthene

22.939min (+0.000) 1.08ng/ul m 08/03/21 JU

response 99373

Ion	Exp%	Act%
252.00	100	100
253.00	21.80	24.26
125.00	12.80	19.59#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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 BNA_P
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Manual Integrations
 APPROVED

mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	411673	20.00	ng/ul	0.00
20) Naphthalene-d8	10.56	136	1731236	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.41	164	982427	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.16	188	1921892	20.00	ng/ul	0.00
79) Chrysene-d12	21.26	240	1737285	20.00	ng/ul	0.00
88) Perylene-d12	23.61	264	1526644	20.00	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.30	96	33464	3.03	ng/uL	0.00
4) Pyridine-d5	3.70	84	389125	11.89	ng/ul	0.00
7) Phenol-d5	6.95	99	598498	15.48	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.12	67	386467	16.06	ng/ul	0.00
11) 2-Chlorophenol-d4	7.31	132	470247	16.37	ng/ul	0.00
15) 4-Methylphenol-d8	8.49	113	428941	14.11	ng/ul	0.00
21) Nitrobenzene-d5	8.93	128	229170	17.78	ng/ul	0.00
24) 2-Nitrophenol-d4	9.65	143	236686	19.23	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.19	165	394743	16.37	ng/ul	0.00
31) 4-Chloroaniline-d4	10.70	131	527879	13.29	ng/ul	0.00
46) Dimethylphthalate-d6	13.83	166	1251223	17.89	ng/ul	0.00
49) Acenaphthylene-d8	14.10	160	1567160	18.24	ng/ul	0.00
54) 4-Nitrophenol-d4	14.61	143	195514	14.08	ng/ul	0.00
60) Fluorene-d10	15.40	176	1047622	17.82	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.52	200	136300	12.72	ng/ul	0.00
73) Anthracene-d10	17.26	188	1530623	17.64	ng/ul	0.00
81) Pyrene-d10	19.50	212	1722488	19.28	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.46	264	1351605	17.42	ng/ul	0.00
Target Compounds						
72) Phenanthrene	17.20	178	296741	2.875	ng/ul	99
80) Fluoranthene	19.18	202	518637	4.710	ng/ul	97
82) Pyrene	19.53	202	488963	4.280	ng/ul	99
85) Benzo(a)anthracene	21.25	228	244652	2.273	ng/ul	96
87) Chrysene	21.30	228	293484	2.844	ng/ul	98
90) Benzo(b)fluoranthene	22.90	252	297717	3.070	ng/ul	96
91) Benzo(k)fluoranthene	22.94	252	99373m>	1.080	ng/ul>	96
93) Benzo(a)pyrene	23.50	252	154370	1.821	ng/ul#	96
94) Indeno(1,2,3-cd)pyrene	26.03	276	108875	1.003	ng/ul	94

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