

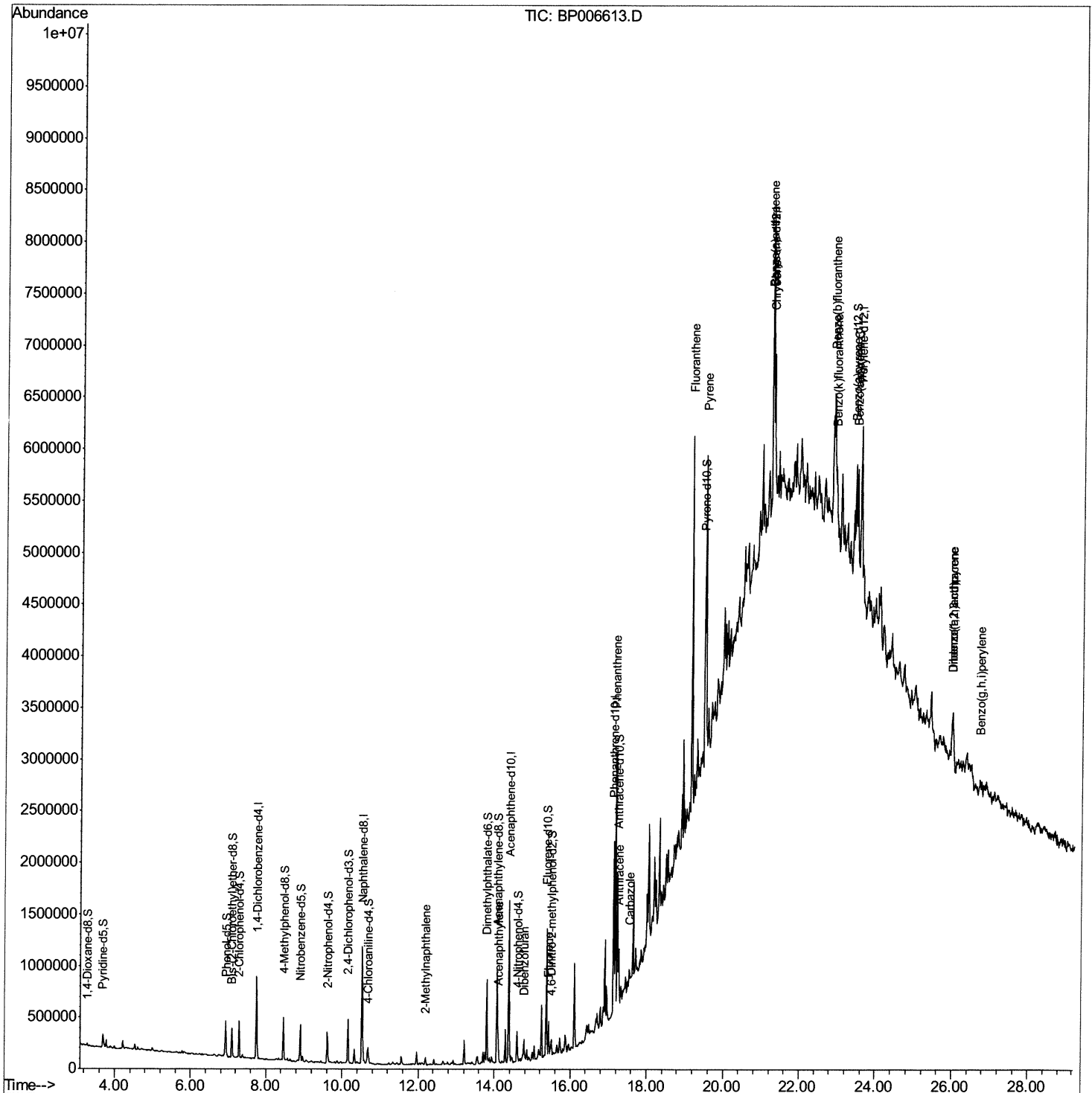
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080921\
Data File : BP006613.D
Acq On : 09 Aug 2021 15:01
Operator : CG/JU
Sample : M3169-21
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00C0

Manual Integrations
APPROVED

mohammad
8/11/2021 9:24:11 PM

Quant Time: Aug 09 16:43:16 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 09 16:04:34 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

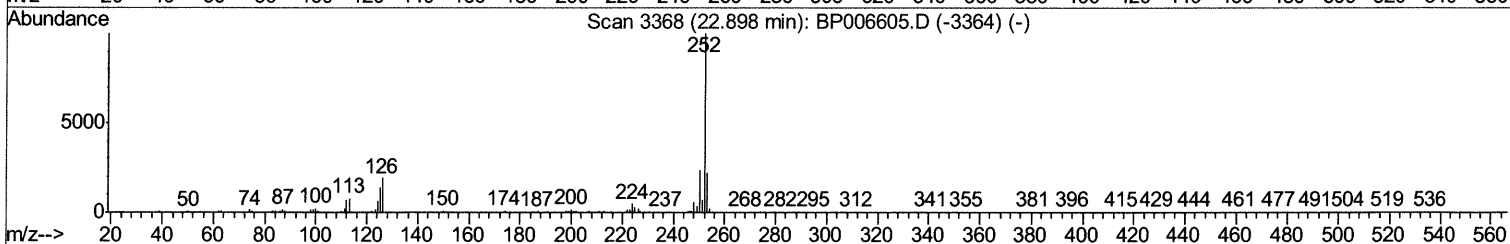
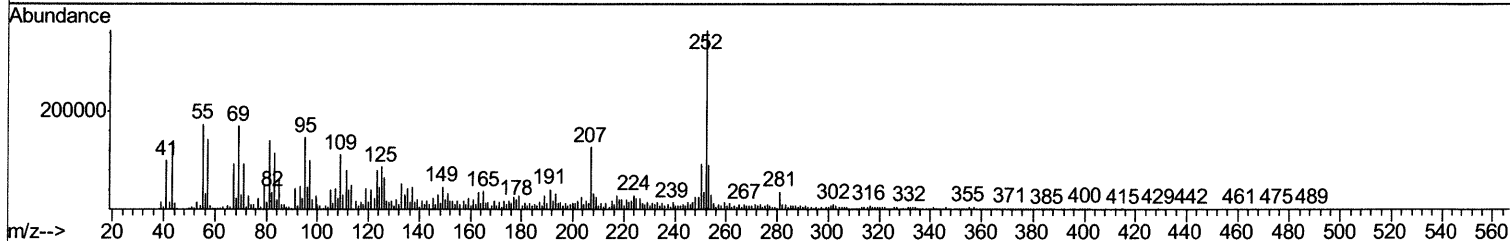
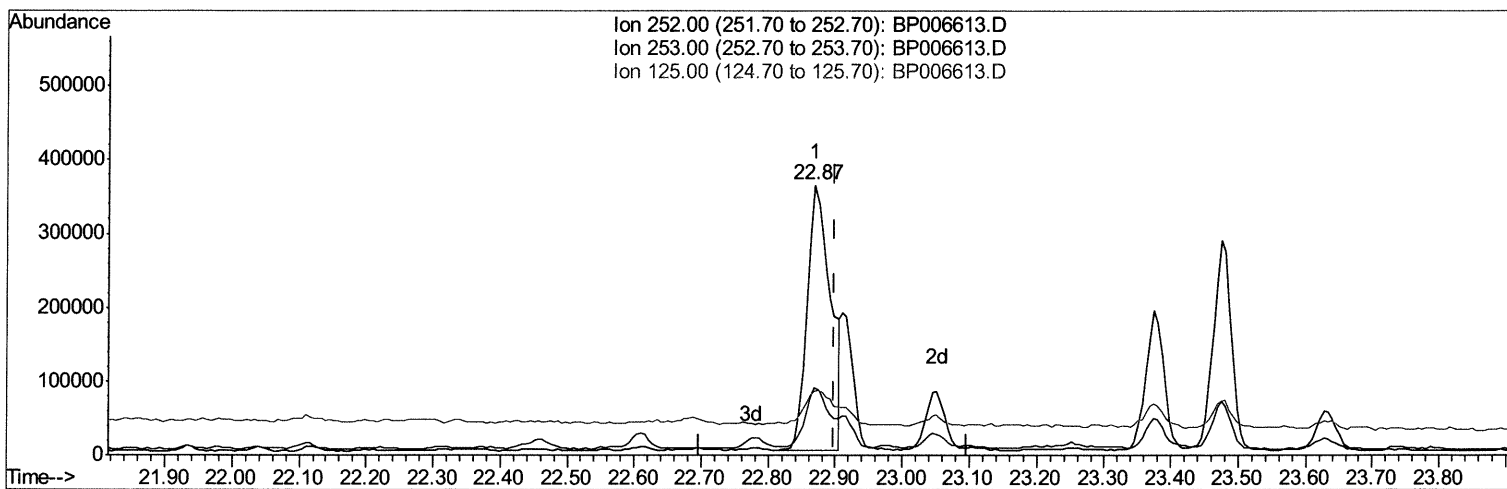
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TIC: BP006613.D

(91) Benzo(k)fluoranthene

22.869min (-0.029) 15.49ng/ul

response 866704

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	24.81
125.00	12.00	23.85#
0.00	0.00	0.00

Quantitation Report (Qedit)

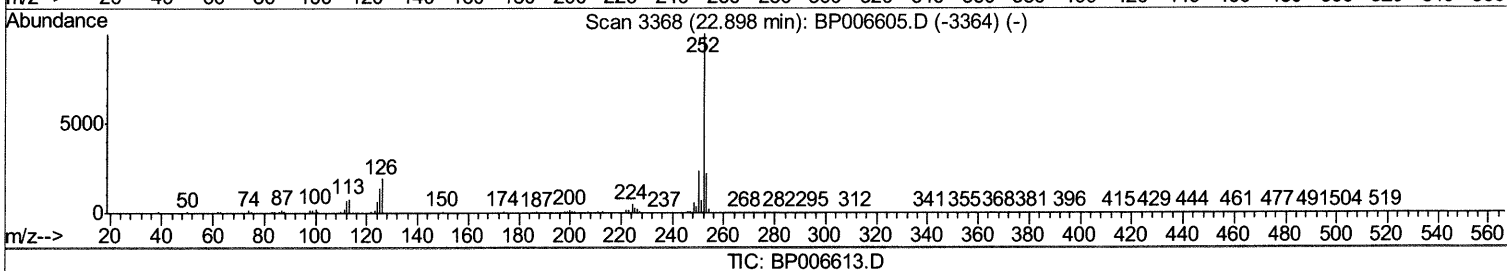
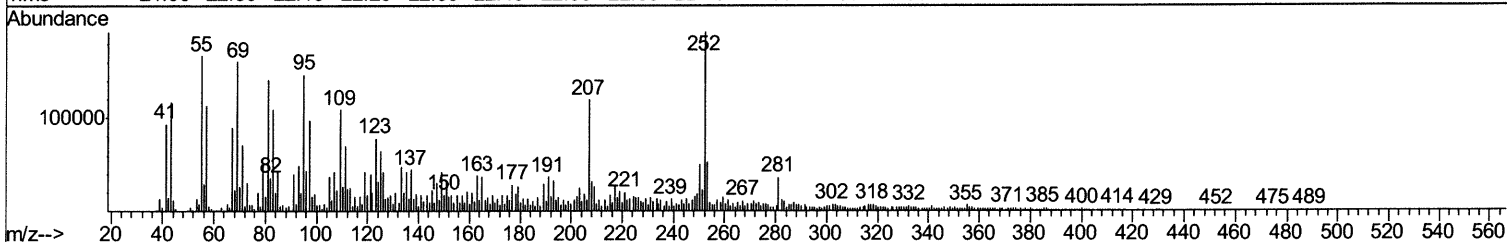
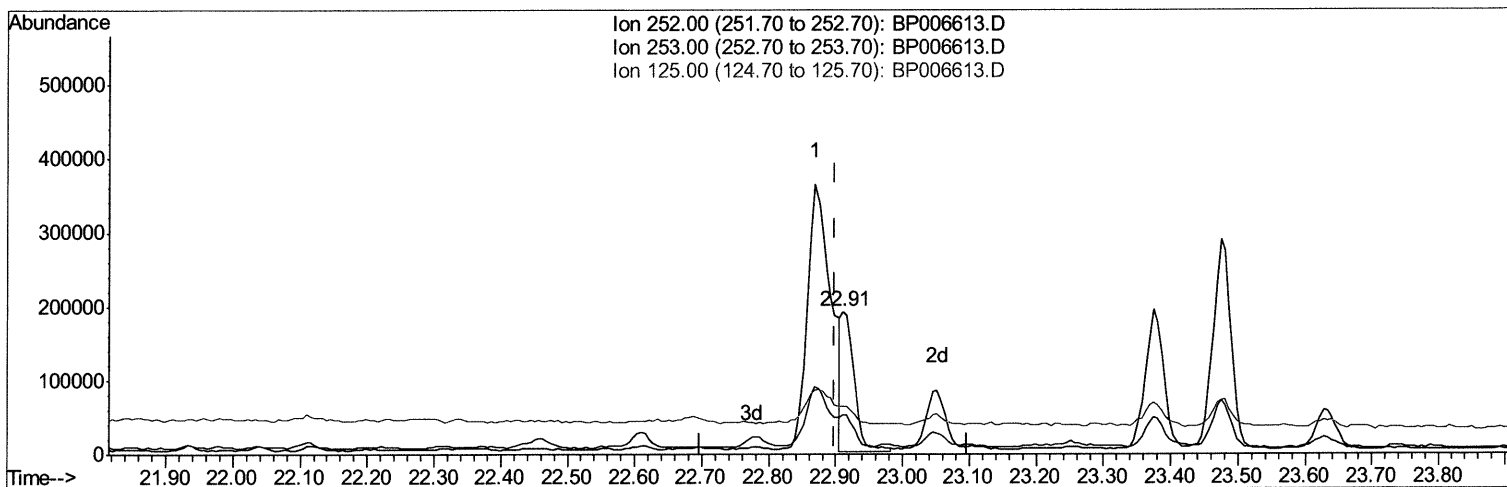
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Response via : Initial Calibration



(91) Benzo(k)fluoranthene

22.910min (+0.012) 4.55ng/ul m 08/12/21 JU

response 254572

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	27.27#
125.00	12.00	33.82#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : M3169-21
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual Integrations
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 09 16:04:34 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	201247	20.00	ng/ul	0.00
20) Naphthalene-d8	10.53	136	873616	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.38	164	481364	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.13	188	904088	20.00	ng/ul	0.00
79) Chrysene-d12	21.24	240	859566	20.00	ng/ul	0.00
88) Perylene-d12	23.58	264	928122	20.00	ng/ul	0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	9702	1.80	ng/uL	0.00
4) Pyridine-d5	3.69	84	90509	5.65	ng/ul	0.01
7) Phenol-d5	6.93	99	197526	10.45	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.09	67	129154	10.98	ng/ul	0.00
11) 2-Chlorophenol-d4	7.29	132	159029	11.33	ng/ul	0.00
15) 4-Methylphenol-d8	8.46	113	152939	10.29	ng/ul	0.00
21) Nitrobenzene-d5	8.90	128	81568	12.54	ng/ul	0.00
24) 2-Nitrophenol-d4	9.62	143	84979	13.68	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.16	165	143826	11.82	ng/ul	0.00
31) 4-Chloroaniline-d4	10.68	131	98350	4.91	ng/ul	0.00
46) Dimethylphthalate-d6	13.80	166	485116	14.16	ng/ul	0.00
49) Acenaphthylene-d8	14.07	160	601419	14.29	ng/ul	0.00
54) 4-Nitrophenol-d4	14.60	143	57185	8.41	ng/ul	0.01
60) Fluorene-d10	15.37	176	419474	14.56	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.50	200	26110	5.18	ng/ul	0.00
73) Anthracene-d10	17.23	188	614993	15.07	ng/ul	0.00
81) Pyrene-d10	19.48	212	676149	15.29	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.43	264	666871	14.14	ng/ul	0.02

Target Compounds

					Ovalue	
36) 2-Methylnaphthalene	12.20	142	33188	1.069	ng/ul	100
50) Acenaphthylene	14.10	152	207130	5.053	ng/ul	99
56) Dibenzofuran	14.78	168	60498	1.500	ng/ul	95
61) Fluorene	15.43	166	133468	4.023	ng/ul#	98
72) Phenanthrene	17.18	178	1439448	29.651	ng/ul	99
74) Anthracene	17.27	178	319732	6.495	ng/ul	97
77) Carbazole	17.55	167	79941	1.775	ng/ul#	91
80) Fluoranthene	19.16	202	1818364	33.378	ng/ul	98
82) Pyrene	19.51	202	1355359	23.976	ng/ul	97
85) Benzo(a)anthracene	21.23	228	778088	14.608	ng/ul#	88
87) Chrysene	21.27	228	729102	14.281	ng/ul#	94
90) Benzo(b)fluoranthene	22.87	252	866704	14.702	ng/ul#	86
91) Benzo(k)fluoranthene	22.91	252	254572m>	4.550	ng/ul>	08/12/21
93) Benzo(a)pyrene	23.47	252	527040	10.228	ng/ul#	85
94) Indeno(1,2,3-cd)pyrene	25.99	276	373520	5.663	ng/ul	96
95) Dibenzo(a,h)anthracene	26.00	278	109534	1.961	ng/ul#	53
96) Benzo(q,h,i)perylene	26.74	276	85090	1.561	ng/ul#	68

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