

(QT Reviewed)

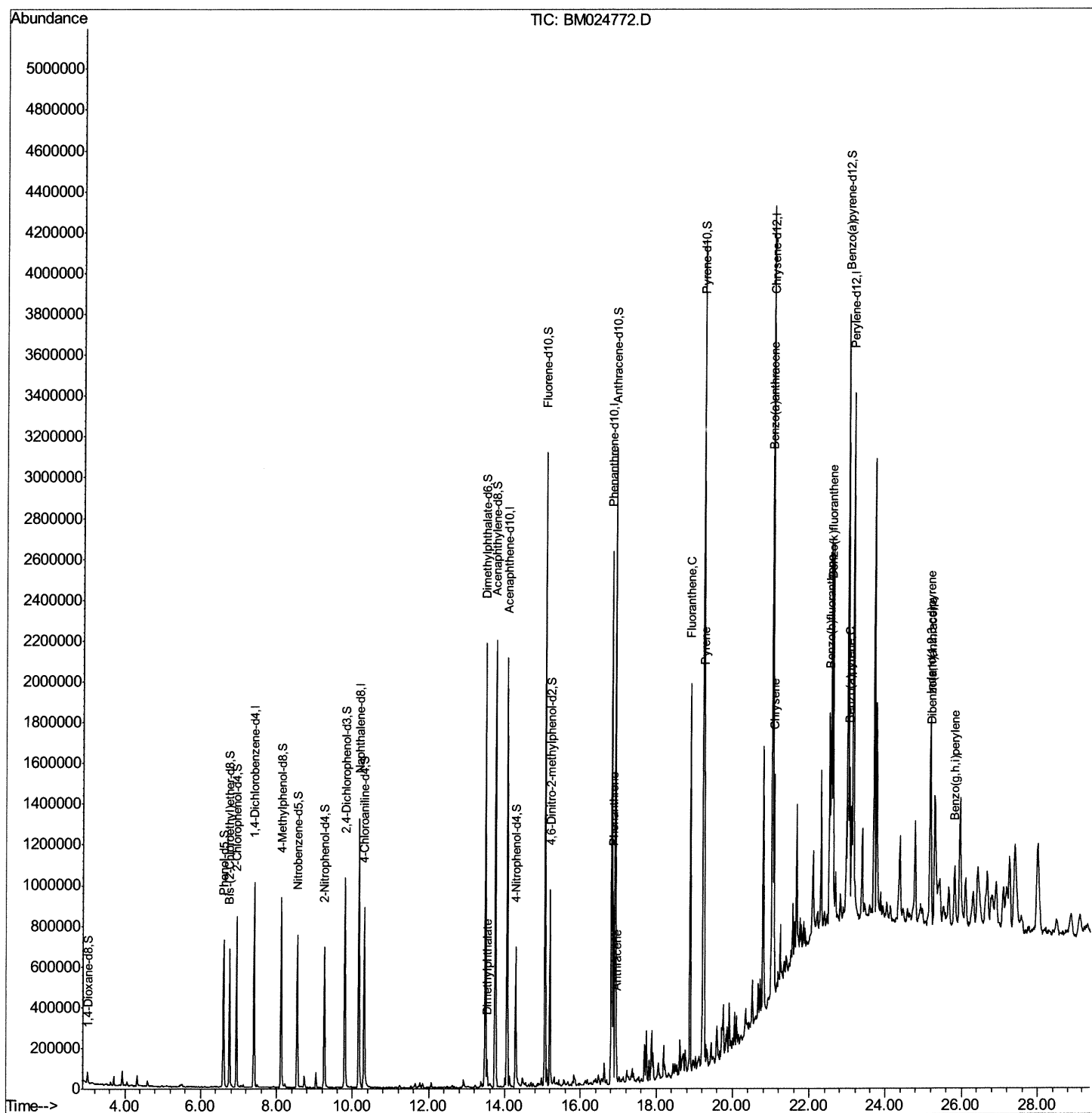
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020620\
Data File : BM024772.D
Acq On : 07 Feb 2020 21:41
Operator : CG/JU
Sample : L1399-17
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB6T2

Manual Integrations
APPROVED

mohammad
2/11/2020 8:22:07 AM

Quant Time: Feb 08 00:52:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM020620MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 07 22:29:17 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

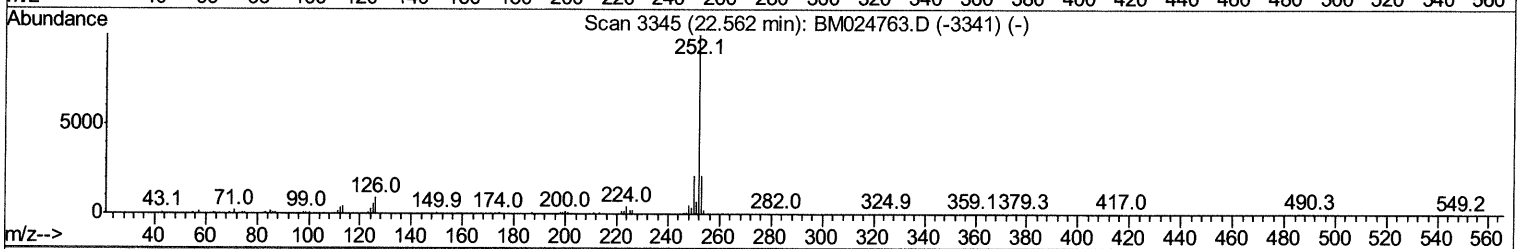
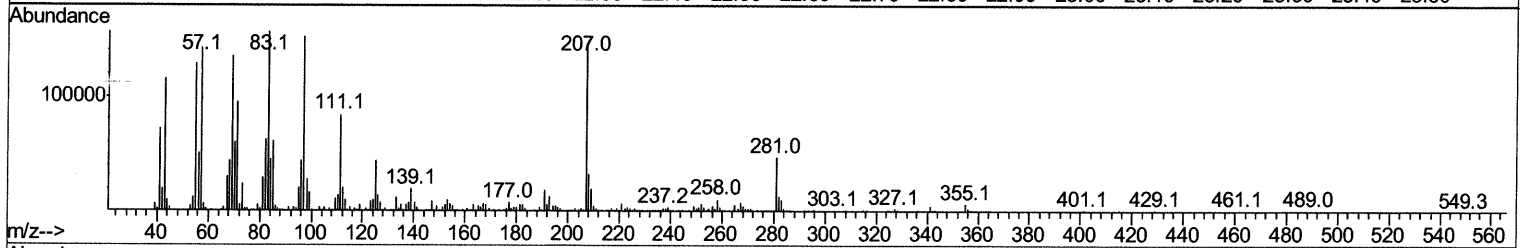
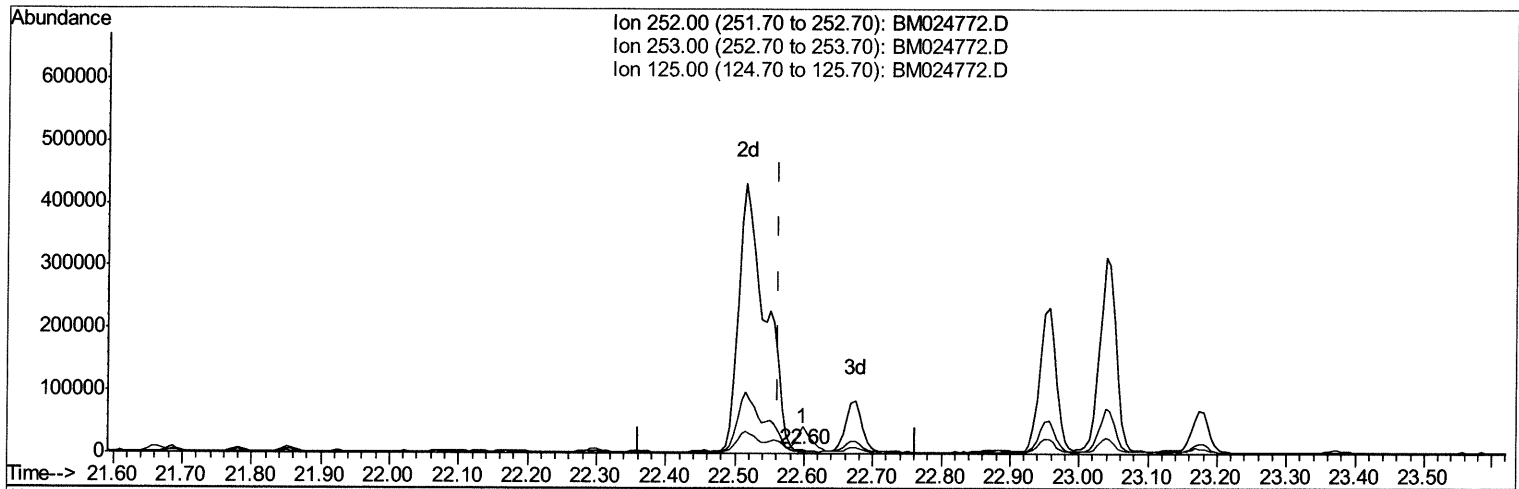
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TIC: BM024772.D

(89) Benzo(k)fluoranthene

22.598min (+0.036) 0.03ng/ul

response 2908

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	56.48#
125.00	6.40	676.93#
0.00	0.00	0.00

Quantitation Report (Qedit)

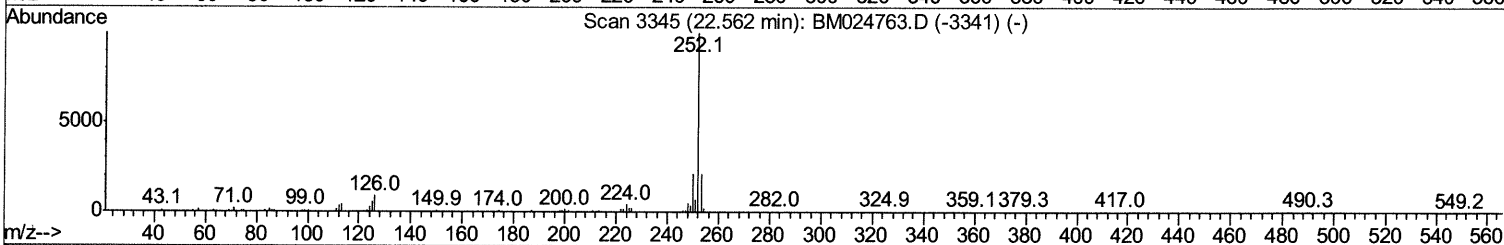
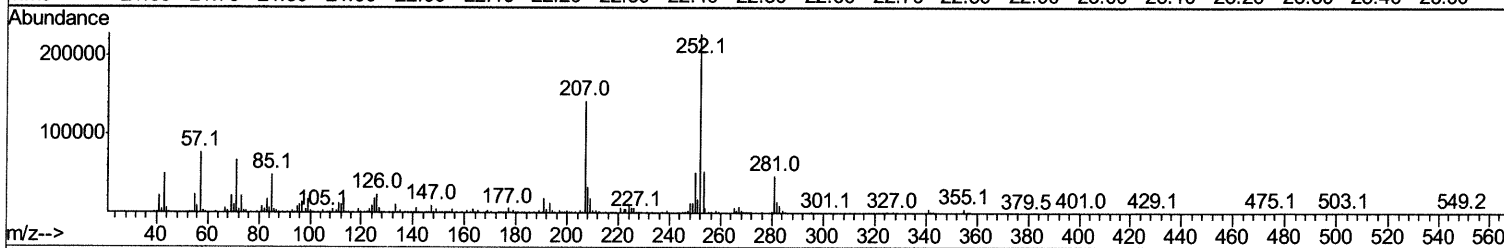
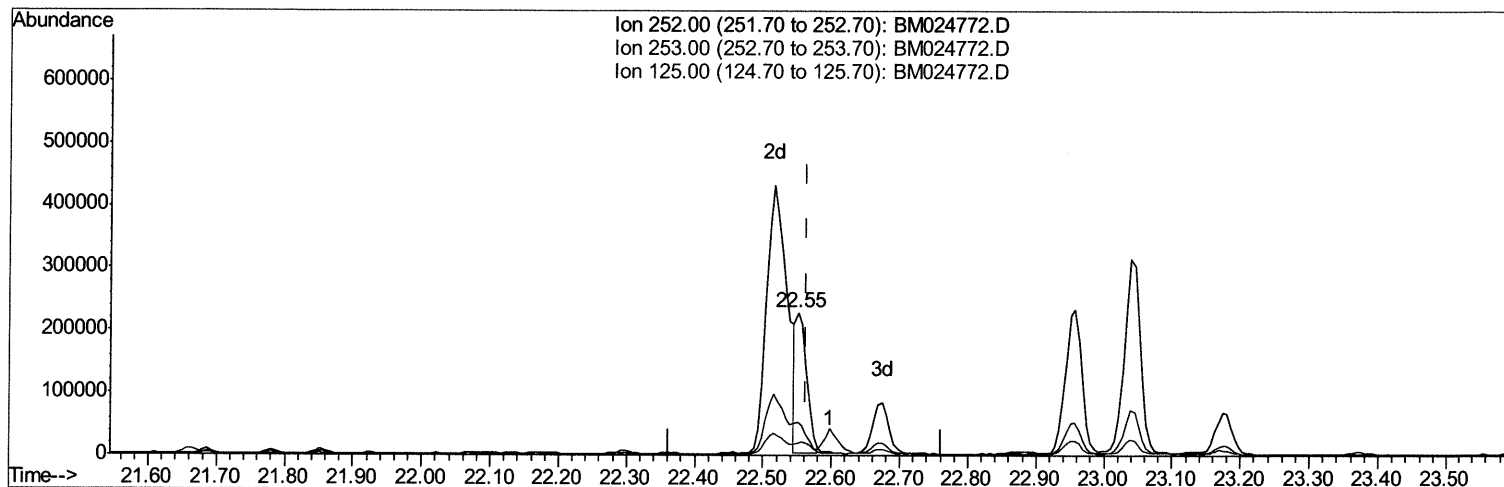
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 Response via : Initial Calibration



TIC: BM024772.D

(89) Benzo(k)fluoranthene

22.551min (-0.011) 2.14ng/ul m JU 02/12/20

response 235409

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	23.22
125.00	6.40	8.30#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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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	7.40	152	272809	20.00	ng/ul	0.00
18) Naphthalene-d8	10.16	136	1082643	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.06	164	733250	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.81	188	1677477	20.00	ng/ul	0.00
78) Chrysene-d12	21.03	240	1724366	20.00	ng/ul	0.00
86) Perylene-d12	23.13	264	1796432	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.02	96	25243	4.45	ng/uL	0.00
5) Phenol-d5	6.60	99	417404	23.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.76	67	269001	26.28	ng/ul	0.00
9) 2-Chlorophenol-d4	6.95	132	380446	23.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.12	113	362981	23.74	ng/ul	-0.01
19) Nitrobenzene-d5	8.55	128	199314	25.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.26	143	242080	26.17	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.80	165	419589	22.49	ng/ul	0.00
29) 4-Chloroaniline-d4	10.30	131	496881	28.47	ng/ul	0.00
44) Dimethylphthalate-d6	13.48	166	1532115	27.68	ng/ul	0.00
47) Acenaphthylene-d8	13.75	160	1564995	24.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.28	143	204775	22.77	ng/ul	0.00
58) Fluorene-d10	15.06	176	1248449	25.47	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	234983	21.49	ng/ul	0.00
71) Anthracene-d10	16.91	188	1811875	23.66	ng/ul	0.00
79) Pyrene-d10	19.22	212	2082373	24.14	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.00	264	2130928	23.62	ng/ul	0.00

Target Compounds

					Ovalue	
45) Dimethylphthalate	13.53	163	88567	1.534	ng/ul	99
70) Phenanthrene	16.85	178	584283	6.388	ng/ul	100
72) Anthracene	16.95	178	124461	1.342	ng/ul	98
77) Fluoranthene	18.88	202	1176236	10.476	ng/ul	99
80) Pyrene	19.24	202	1035542	9.073	ng/ul	98
83) Benzo(a)anthracene	21.01	228	612735	5.275	ng/ul	96
85) Chrysene	21.06	228	624441	5.461	ng/ul	99
88) Benzo(b)fluoranthene	22.52	252	916444	8.023	ng/ul#	97
89) Benzo(k)fluoranthene	22.55	252	235409m	2.141	ng/ul	> 97
91) Benzo(a)pyrene	23.04	252	518513	5.063	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.19	276	408676	3.205	ng/ul	98
93) Dibenzo(a,h)anthracene	25.19	278	119325	1.098	ng/ul#	95
94) Benzo(g,h,i)perylene	25.80	276	332632	3.137	ng/ul	97

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

2020/02/11