

(QT Reviewed)

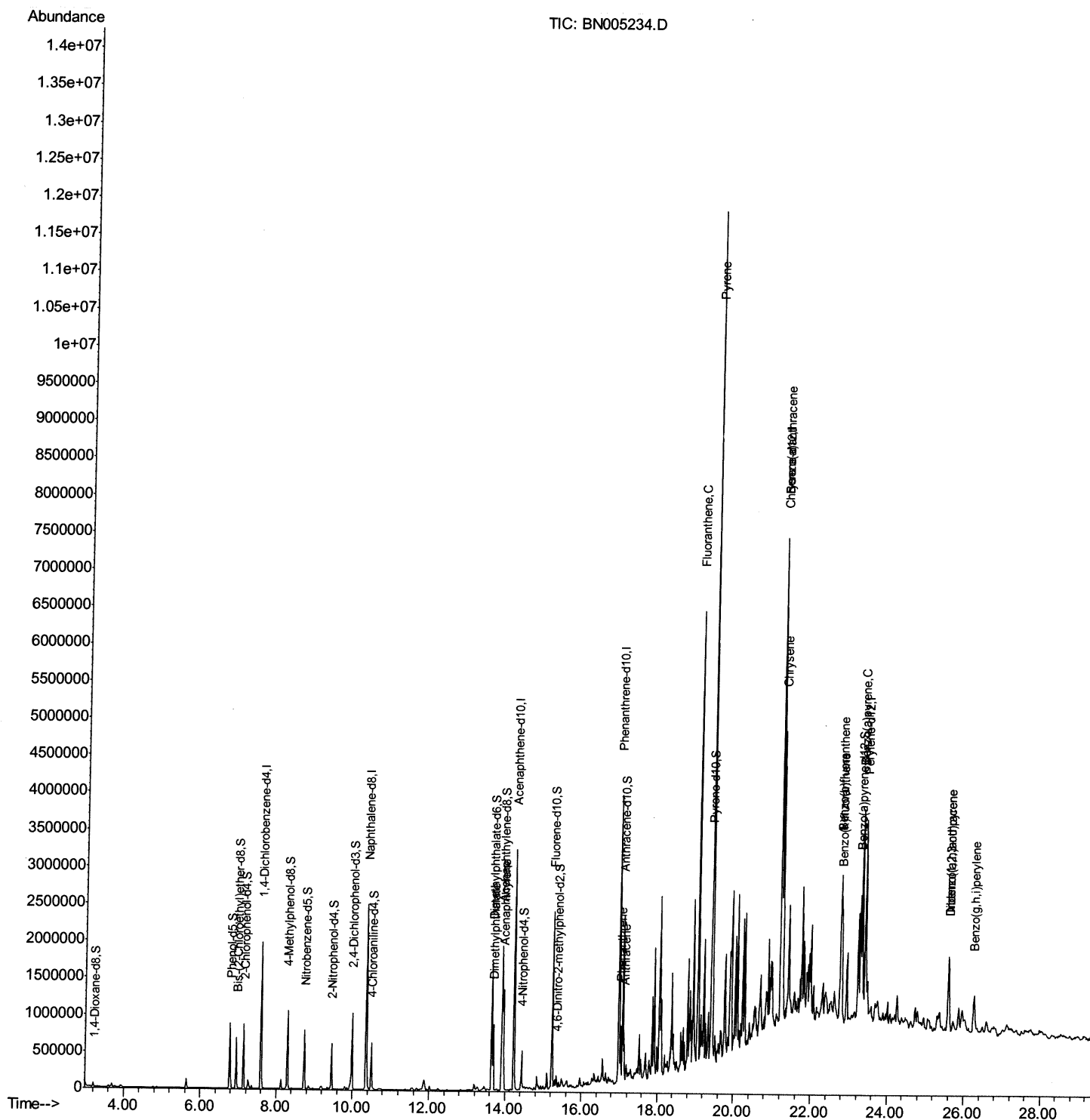
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042319\
Data File : BN005234.D
Acq On : 24 Apr 2019 03:05
Operator : JU/SJ
Sample : K2455-04DL 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GAHQ6DL

Manual Integrations
APPROVED

Sohil
4/24/2019 11:45:57 AM

Quant Time: Apr 24 09:42:13 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 23 03:50:45 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

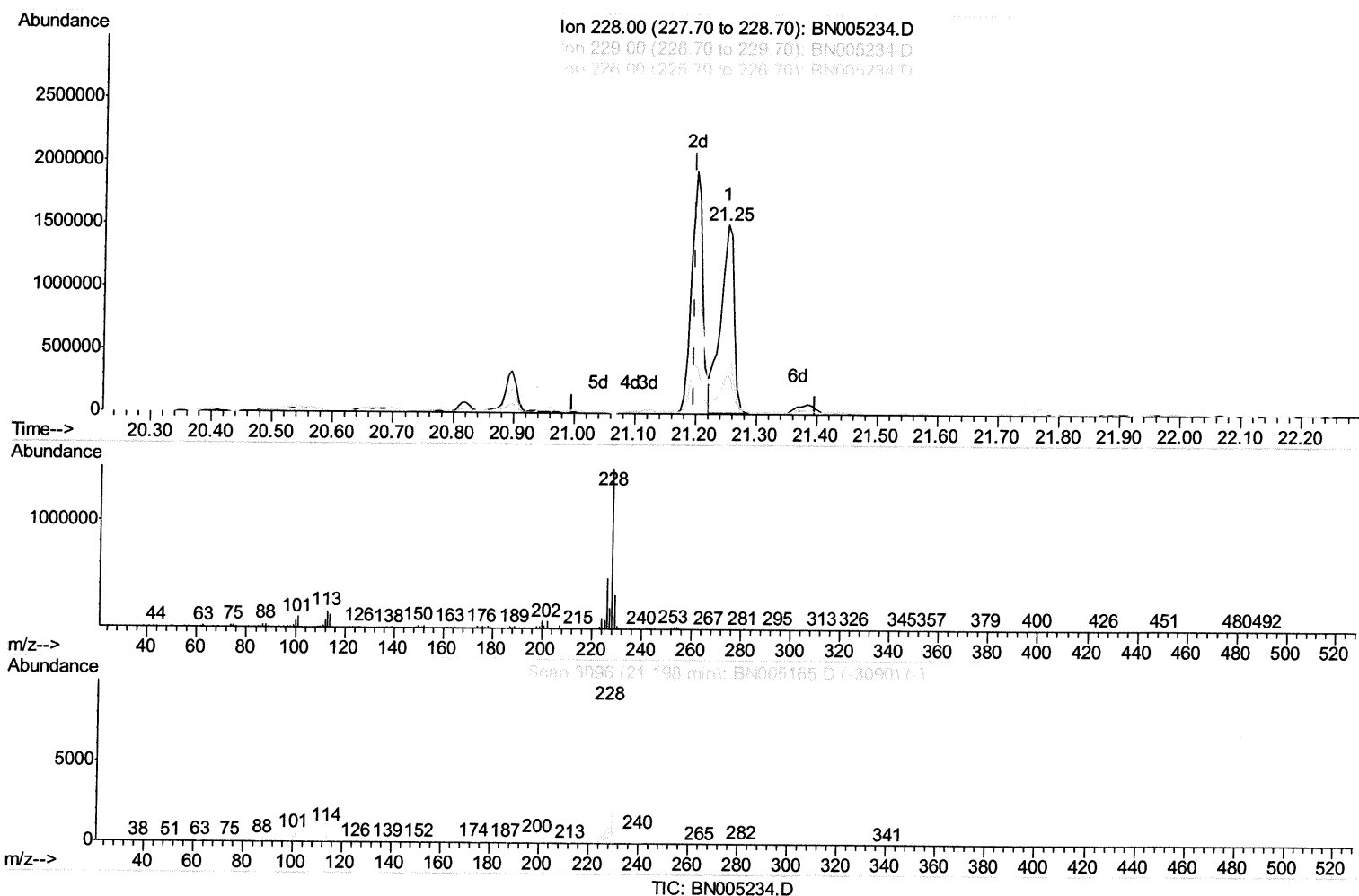
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(82) Benzo(a)anthracene

21.251min (+0.053) 21.85ng/ul

response 2292320

Ion	Exp%	Act%
228.00	100	100
229.00	19.70	20.93
226.00	26.80	31.31
0.00	0.00	0.00

Quantitation Report (Qedit)

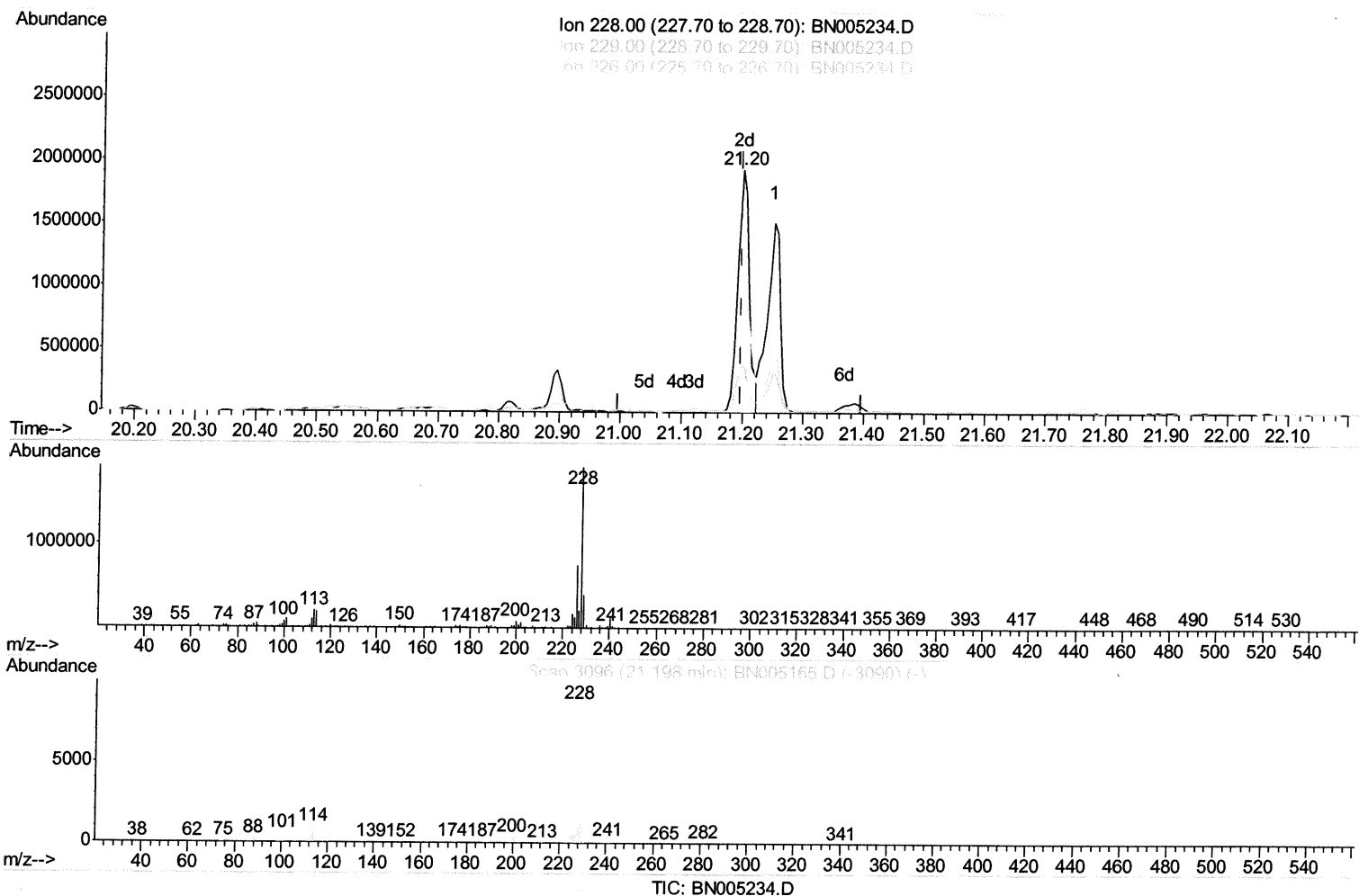
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(82) Benzo(a)anthracene

21.198min (-0.000) 24.13ng/ul mSJ 04/30/19

response 2531435

Ion	Exp%	Act%
228.00	100	100
229.00	19.70	20.53
226.00	26.80	38.84#
0.00	0.00	0.00

Quantitation Report (Qedit)

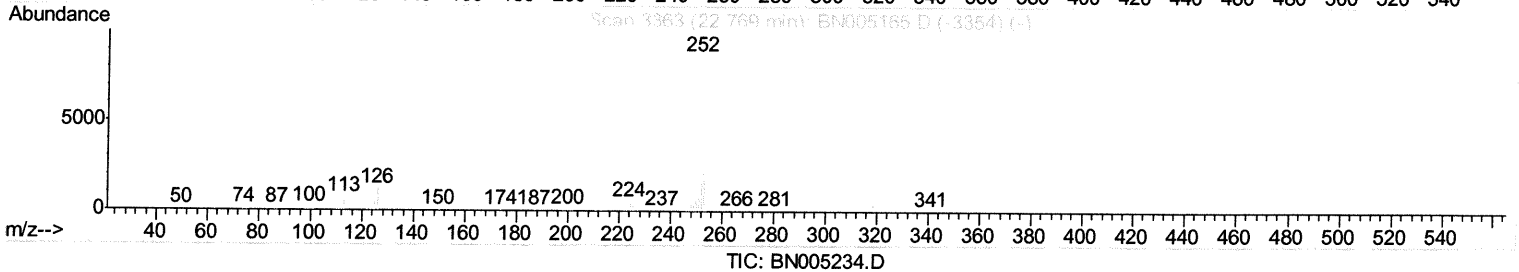
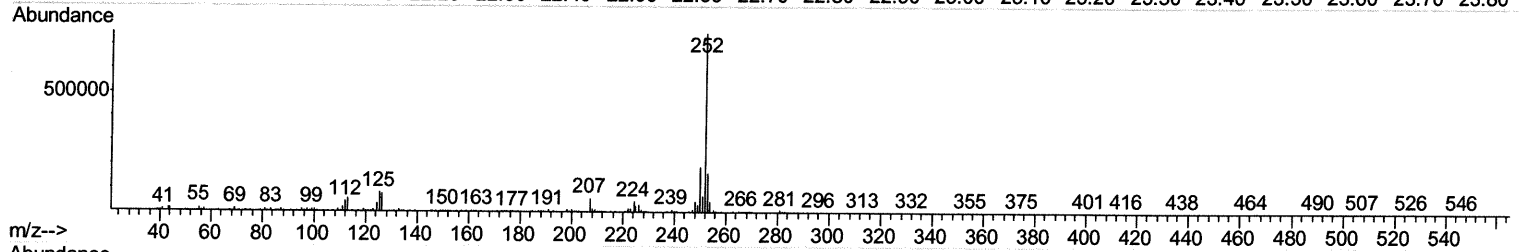
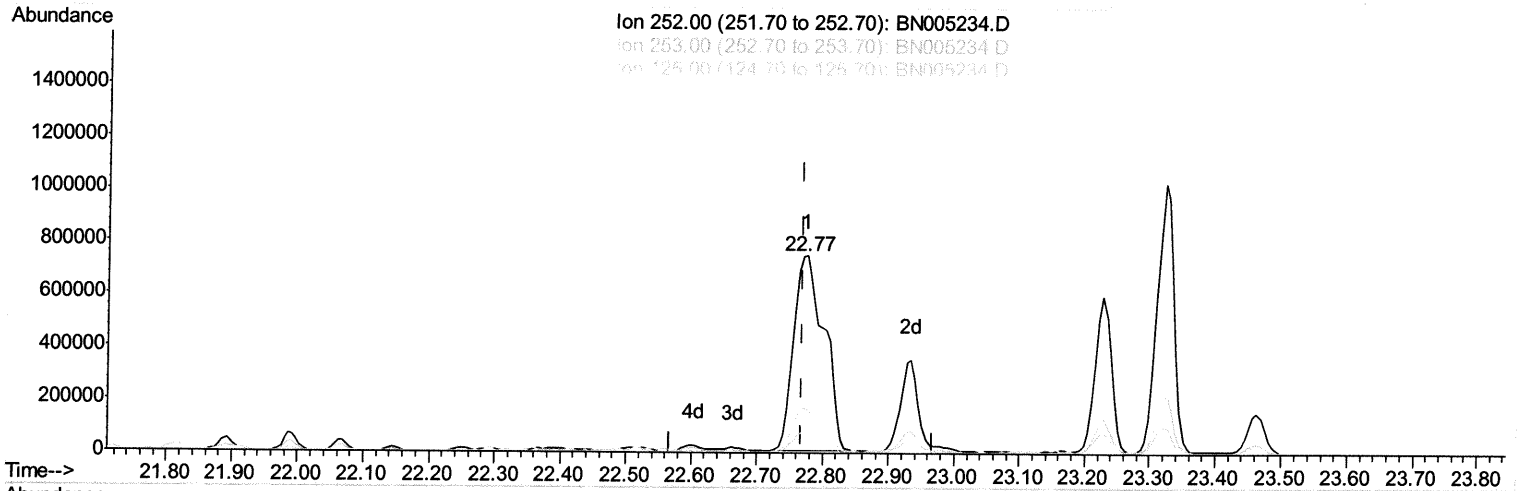
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042319\
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 Acq On : 24 Apr 2019 03:05
 Operator : JU/SJ
 Sample : K2455-04DL 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
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(87) Benzo(b)fluoranthene

22.774min (+0.006) 22.47ng/ul

response 2391782

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	22.10
125.00	10.50	10.76
0.00	0.00	0.00

Quantitation Report (Qedit)

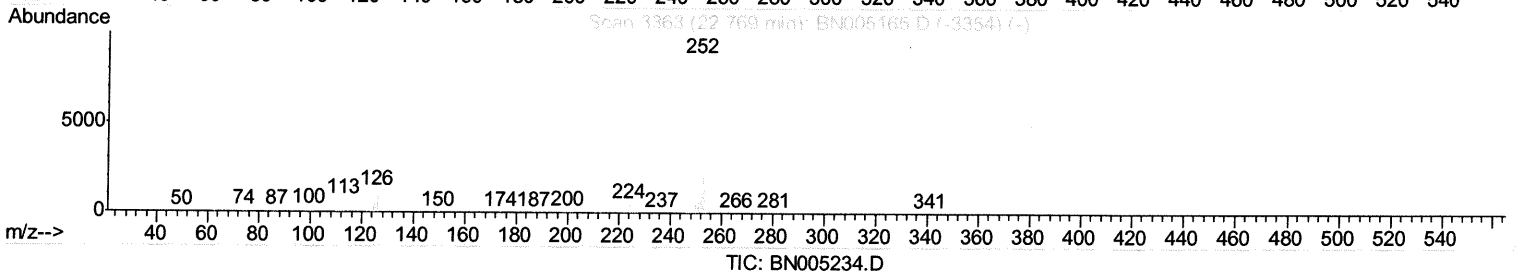
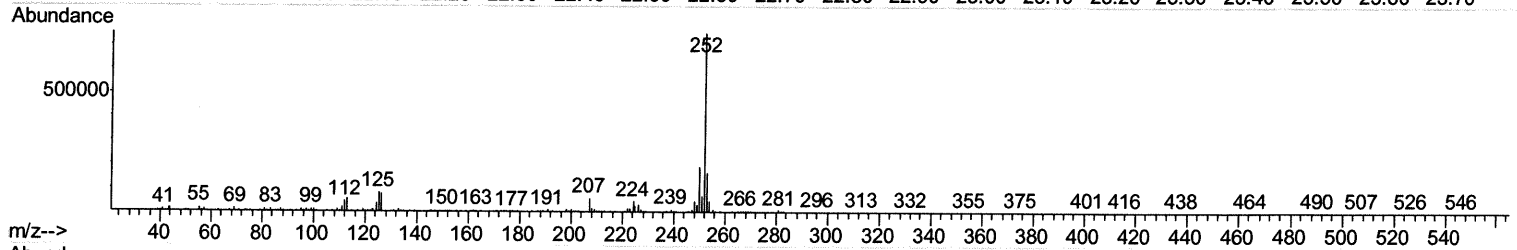
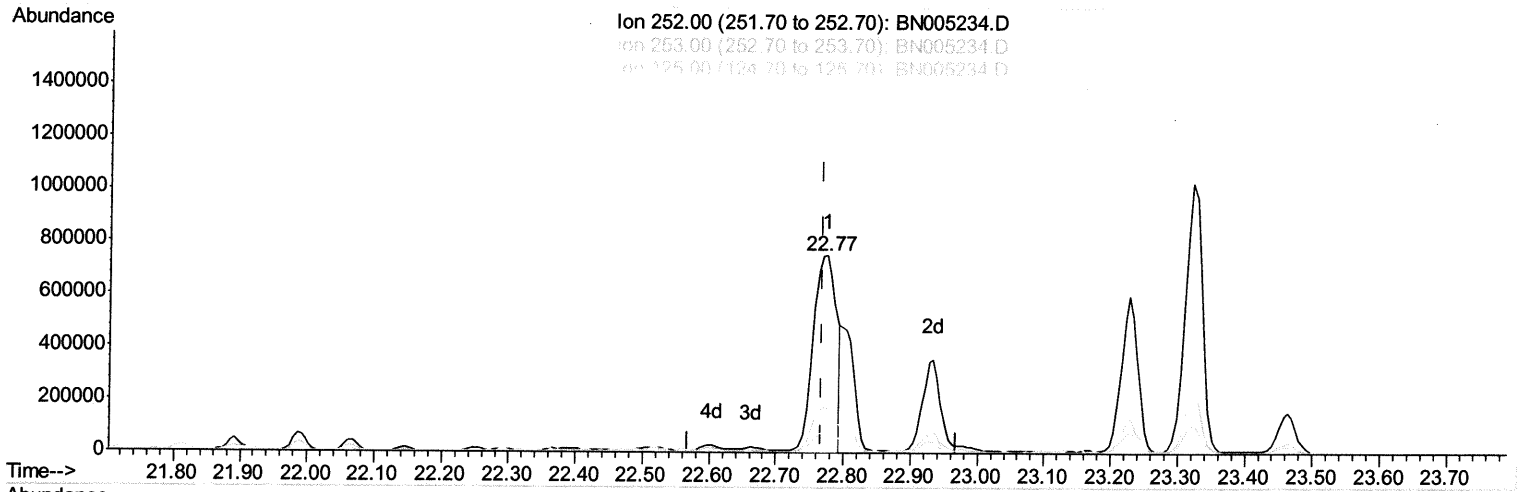
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 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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(87) Benzo(b)fluoranthene

22.774min (+0.006) 16.86ng/ul mSJ 04/30/19

response 1795198

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	22.10
125.00	10.50	10.76
0.00	0.00	0.00

Quantitation Report (Qedit)

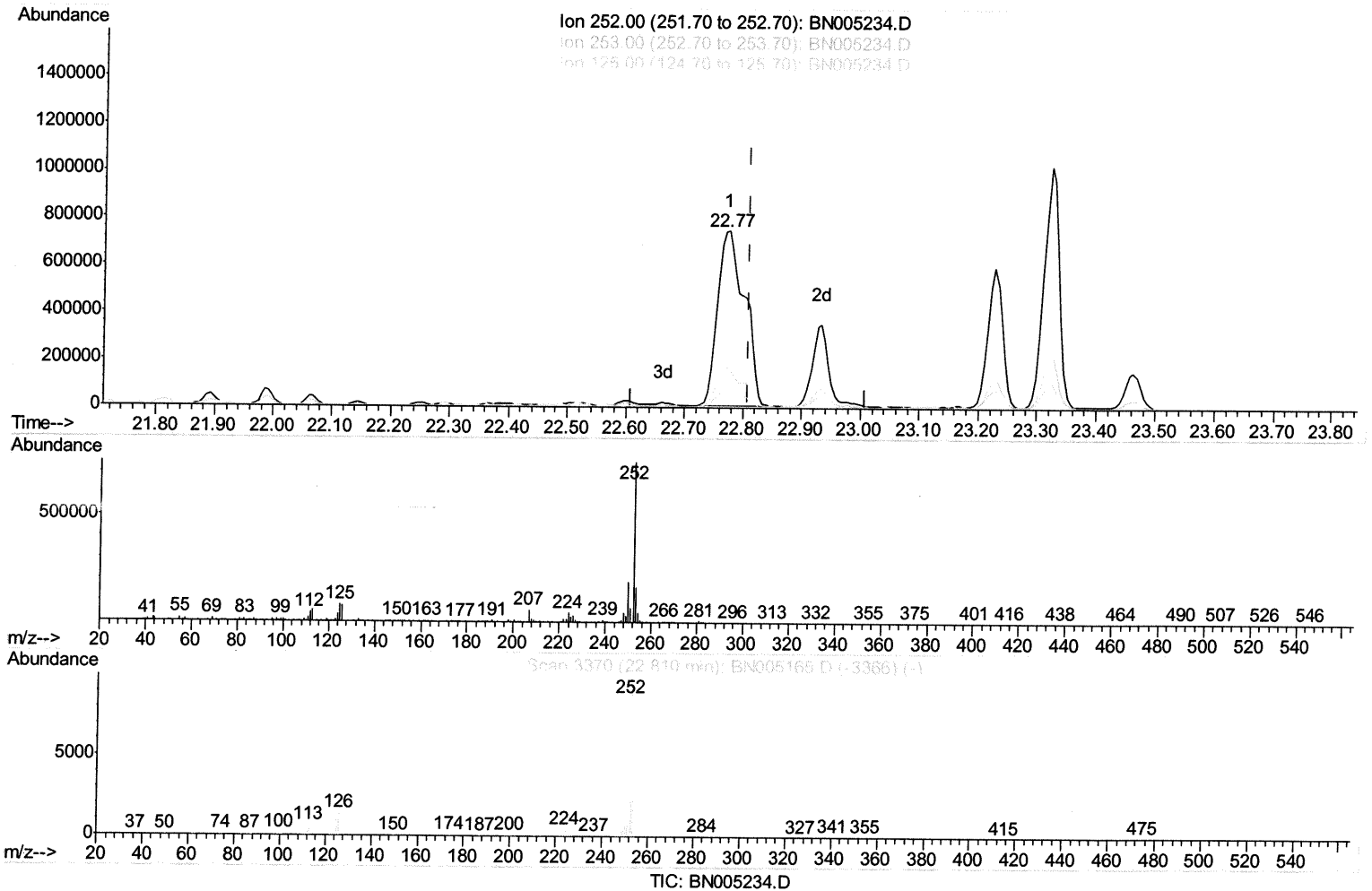
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 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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(88) Benzo(k)fluoranthene
 22.774min (-0.035) 23.52ng/ul
 response 2391782

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	22.10
125.00	9.40	10.76
0.00	0.00	0.00

Quantitation Report (Qedit)

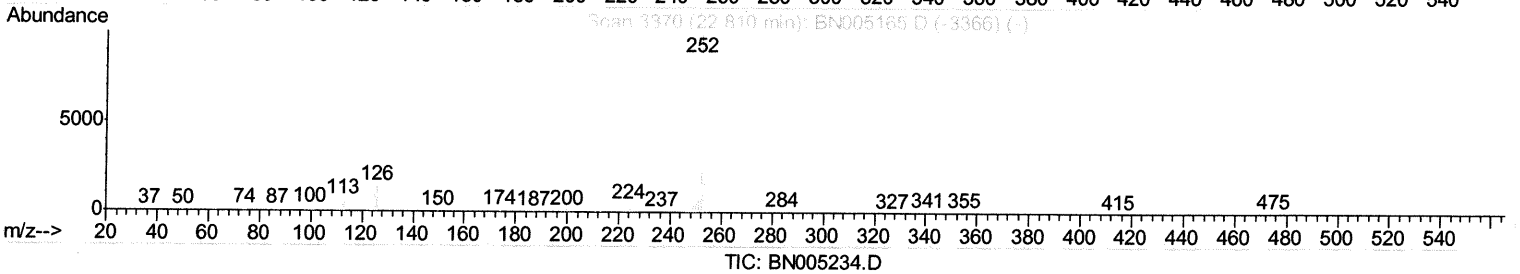
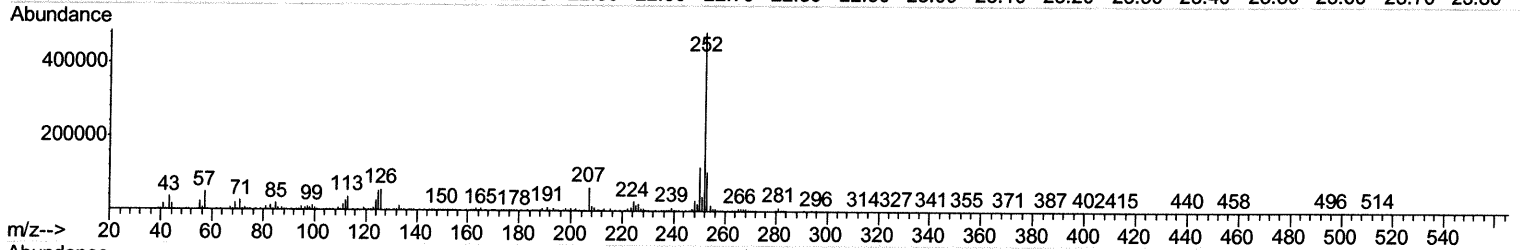
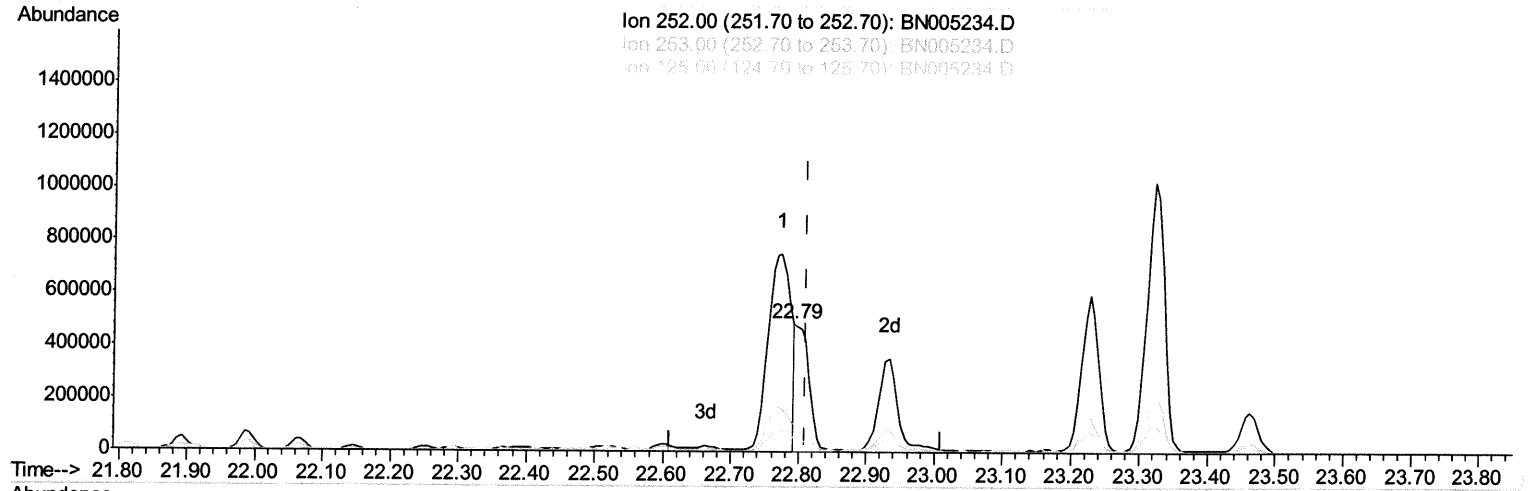
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN042319\
 Data File : BN005234.D
 Acq On : 24 Apr 2019 03:05
 Operator : JU/SJ
 Sample : K2455-04DL 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleID :
 GAHQ6DL

Manual Integrations
 APPROVED

Sohil
 4/24/2019 11:45:57 AM

Quant Time: Apr 24 08:55:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 23 03:50:45 2019
 Response via : Initial Calibration



(88) Benzo(k)fluoranthene

22.792min (-0.018) 6.52ng/ul m SJ04/30/19

response 662964

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	22.04
125.00	9.40	11.11
0.00	0.00	0.00

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 Operator : JU/SJ
 Sample : K2455-04DL 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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 ClientSampled :
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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	7.60	152	492258	20.00	ng/ul	0.00
18) Naphthalene-d8	10.36	136	1975086	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.23	164	1090123	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.99	188	2162055	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	1686045	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	1893902	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.22	96	20029	1.96	ng/uL	0.00
5) Phenol-d5	6.78	99	462528	11.96	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	283415	11.75	ng/ul	0.00
9) 2-Chlorophenol-d4	7.13	132	363974	12.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	374230	12.38	ng/ul	0.00
19) Nitrobenzene-d5	8.74	128	174913	14.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.45	143	180424	14.64	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	370044	13.10	ng/ul	0.00
29) 4-Chloroaniline-d4	10.50	131	323544	11.11	ng/ul	0.00
43) Dimethylphthalate-d6	13.65	166	1104221	13.95	ng/ul	0.00
46) Acenaphthylene-d8	13.92	160	1352812	13.92	ng/ul	0.00
51) 4-Nitrophenol-d4	14.43	143	120502	9.64	ng/ul	0.00
57) Fluorene-d10	15.23	176	891192	13.42	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.35	200	30976	3.19	ng/ul	0.00
70) Anthracene-d10	17.08	188	1270908	14.02	ng/ul	0.00
78) Pyrene-d10	19.40	212	1258078	14.59	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.28	264	1141417	13.79	ng/ul	0.00
Target Compounds						
44) Dimethylphthalate	13.70	163	354394	4.495	ng/ul	Qvalue 99
47) Acenaphthylene	13.95	152	943672	9.880	ng/ul	99
69) Phenanthrene	17.03	178	463471	4.402	ng/ul	99
71) Anthracene	17.12	178	437874	4.082	ng/ul	99
76) Fluoranthene	19.06	202	3458729	29.614	ng/ul	99
79) Pyrene	19.43	202	6962315	64.134	ng/ul	99
82) Benzo(a)anthracene	21.20	228	2531435m	24.134	ng/ul	→ 97 SJ 04/30/19
84) Chrysene	21.25	228	2305827	23.687	ng/ul	97
87) Benzo(b)fluoranthene	22.77	252	1795198m	16.862	ng/ul	→ 97 SJ 04/30/19
88) Benzo(k)fluoranthene	22.79	252	662964m	6.520	ng/ul	→ 97 SJ 04/30/19
90) Benzo(a)pyrene	23.32	252	1809888	18.086	ng/ul	99
91) Indeno(1,2,3-cd)pyrene	25.59	276	816567	7.381	ng/ul	97
92) Dibenzo(a,h)anthracene	25.60	278	259955	2.769	ng/ul#	93
93) Benzo(g,h,i)perylene	26.25	276	592284	6.474	ng/ul	99

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