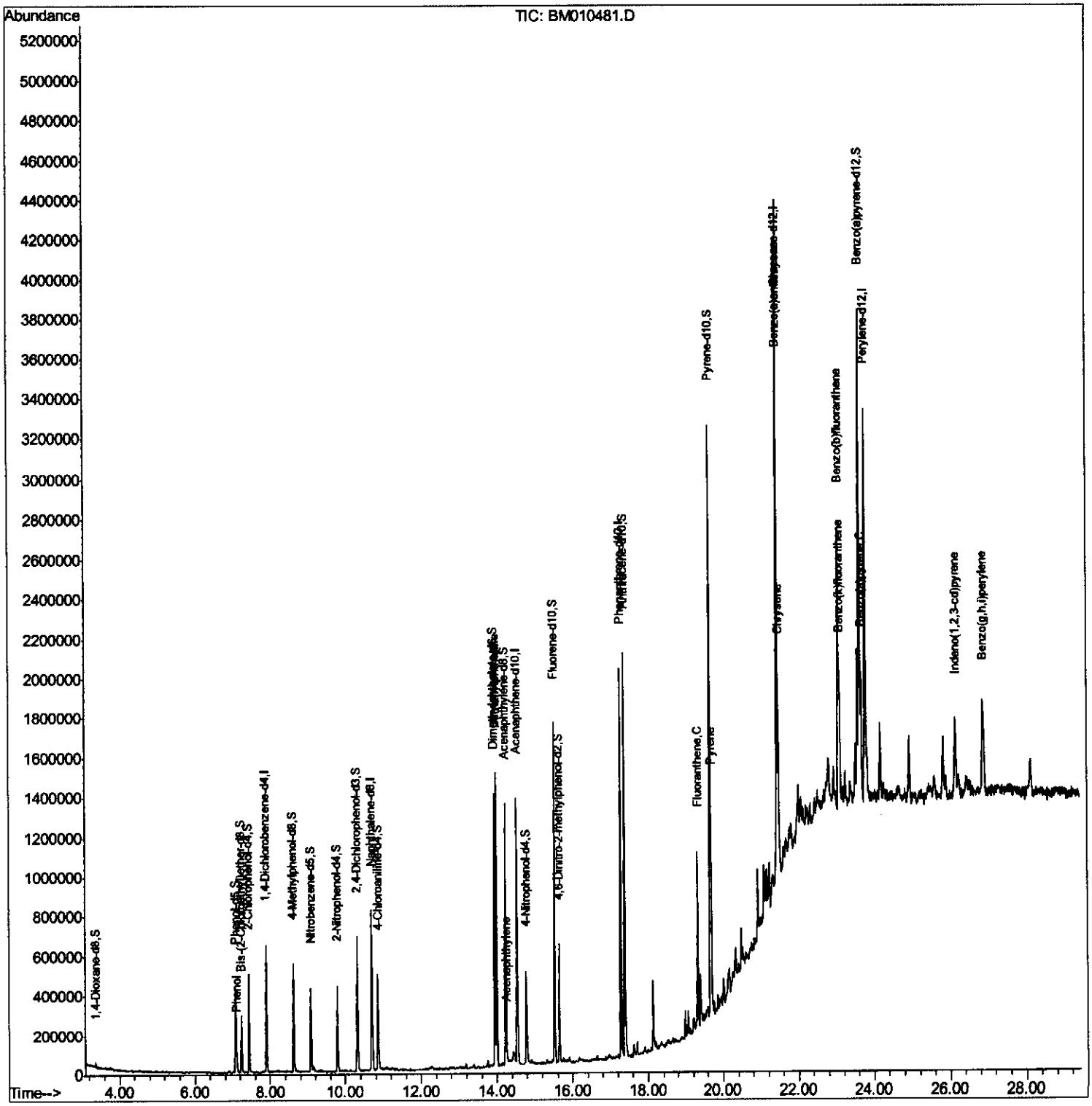
Quantitation Report (LSC Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061017\
Data File : BM010481.D
Acq On : 10 Jun 2017 12:34
Operator : SJ/MA
Sample : I3521-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jun 12 06:04:48 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 12 04:15:09 2017
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
F5C48

Manual Integrations APPROVED
mohammad
6/12/2017 3:40:32 PM



Quantitation Report (Qedit)

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Data File : BM010481.D
Acq On : 10 Jun 2017 12:34
Operator : SJ/MA
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Quant Title : SVOA CALIBRATION
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Instrument :
BNA_M
ClientSampleId :
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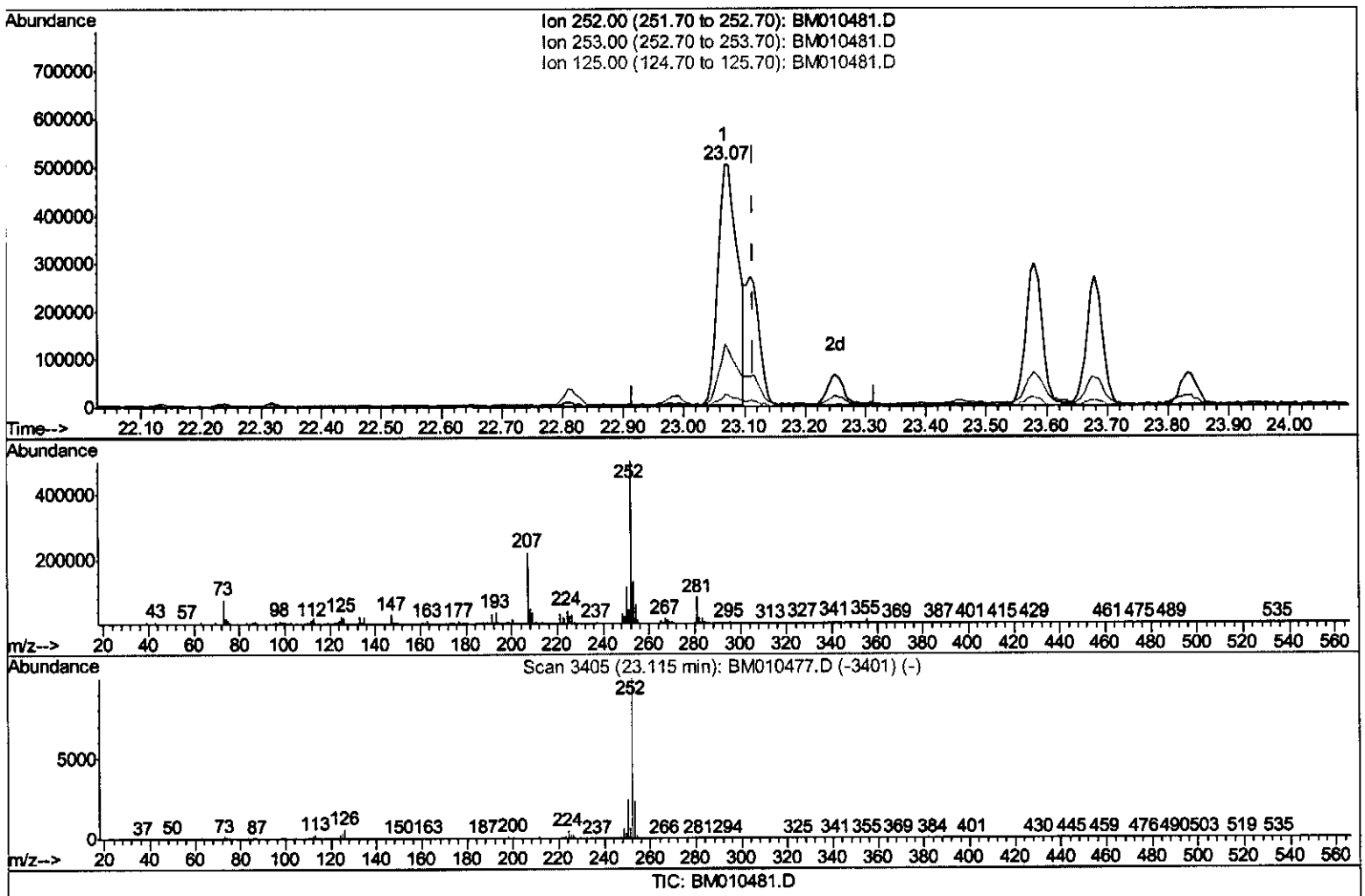
Manual Integrations APPROVED
mohammad
6/12/2017 3:40:32 PM



TIC: BM010481.D

(86) Benzo(k)fluoranthene

23.068min (-0.047) 13.66ng/ul

response 1129455

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	25.80#
125.00	4.30	5.18#
0.00	0.00	0.00

Quantitation Report (Qedit)

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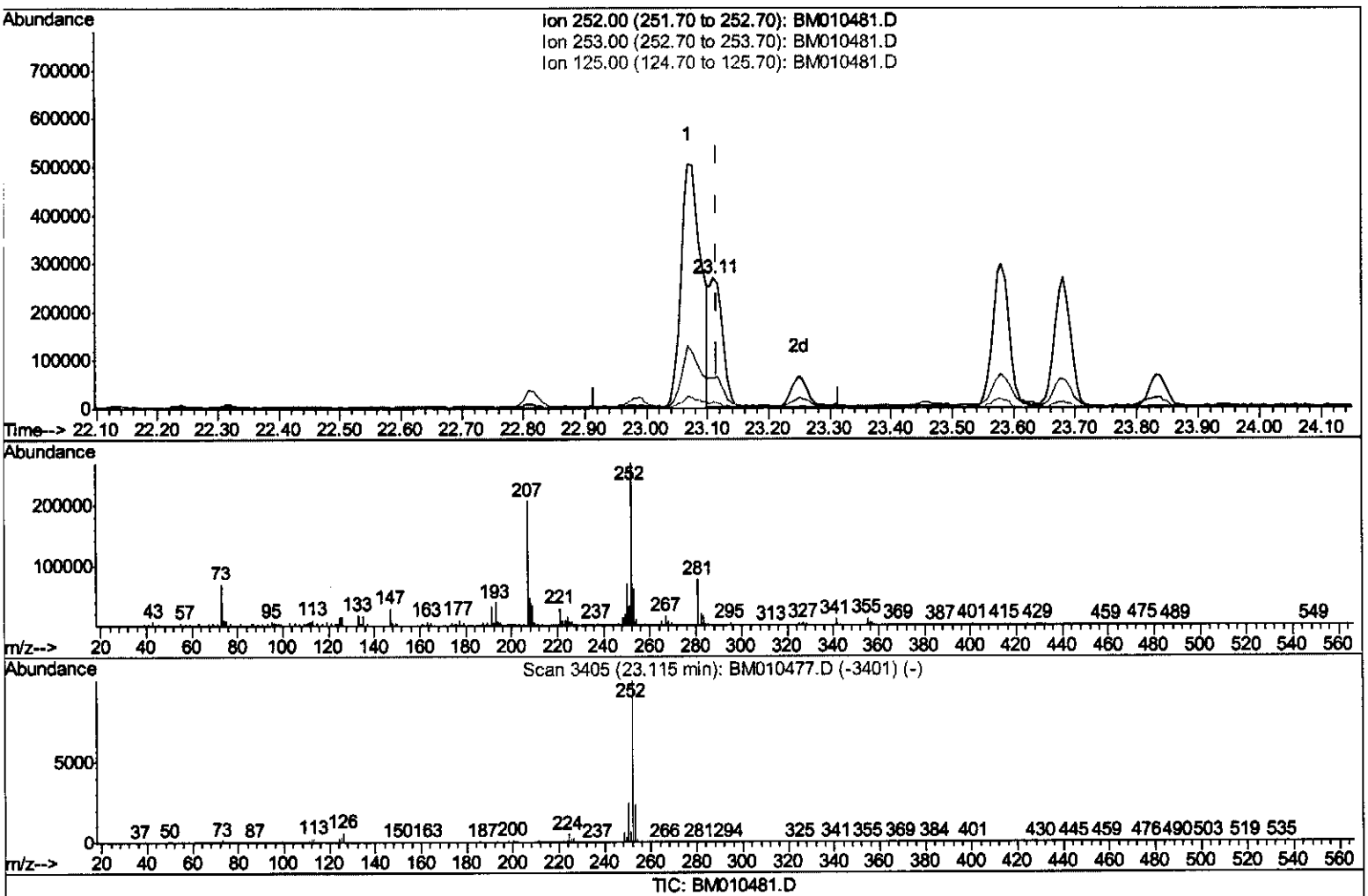
Manual Integrations APPROVED
mohammad
6/12/2017 3:40:32 PM



TIC: BM010481.D

(86) Benzo(k)fluoranthene

23.109min (-0.006) 5.08ng/ul m SJ 6/19/17

response 420064

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	22.79
125.00	4.30	5.42#
0.00	0.00	0.00

Quantitation Report (LSC Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061017\
Data File : BM010481.D
Acq On : 10 Jun 2017 12:34
Operator : SJ/MA
Sample : I3521-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F5C48

Manual Integrations APPROVED
mohammad
6/12/2017 3:40:32 PM

Quant Time: Jun 12 06:04:48 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 12 04:15:09 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	151966	20.00	ng/ul	0.00
18) Naphthalene-d8	10.70	136	554635	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.53	164	383698	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.28	188	977189	20.00	ng/ul	0.00
75) Chrysene-d12	21.44	240	1431382	20.00	ng/ul	0.00
83) Perylene-d12	23.79	264	1487006	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.36	96	12086	3.26	ng/uL	0.00
5) Phenol-d5	7.08	99	193939	16.84	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.23	67	117004	18.08	ng/ul	0.00
9) 2-Chlorophenol-d4	7.43	132	170404	18.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.62	113	168077	18.09	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	81400	20.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.79	143	102139	21.70	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	216246	22.37	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	200890	21.54	ng/ul	0.00
43) Dimethylphthalate-d6	13.94	166	738579	23.17	ng/ul	0.00
46) Acenaphthylene-d8	14.23	160	796081	21.41	ng/ul	0.00
51) 4-Nitrophenol-d4	14.77	143	89009	18.66	ng/ul	0.00
57) Fluorene-d10	15.53	176	638448	22.78	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.65	200	130432	18.84	ng/ul	0.00
70) Anthracene-d10	17.38	188	1088666	22.90	ng/ul	0.00
76) Pyrene-d10	19.67	212	1424800	24.07	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.63	264	1631616	23.57	ng/ul	0.00
Target Compounds						Qvalue
6) Phenol	7.11	94	16609	1.407	ng/ul#	82
44) Dimethylphthalate	13.99	163	791623	25.249	ng/ul	99
47) Acenaphthylene	14.26	152	36689	1.010	ng/ul	98
74) Fluoranthene	19.33	202	518135	8.077	ng/ul#	96
77) Pyrene	19.70	202	605660	8.211	ng/ul#	93
80) Benzo(a)anthracene	21.43	228	416536	5.160	ng/ul	96
82) Chrysene	21.48	228	594004	8.141	ng/ul	95
85) Benzo(b)fluoranthene	23.07	252	1129455	13.053	ng/ul	94
86) Benzo(k)fluoranthene	23.11	252	420064m	5.082	ng/ul	
88) Benzo(a)pyrene	23.68	252	489125	5.703	ng/ul	99
89) Indeno(1,2,3-cd)pyrene	26.16	276	422405	3.898	ng/ul	96
91) Benzo(g,h,i)perylene	26.90	276	330101	3.698	ng/ul	98

SJ 06/19/17

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