

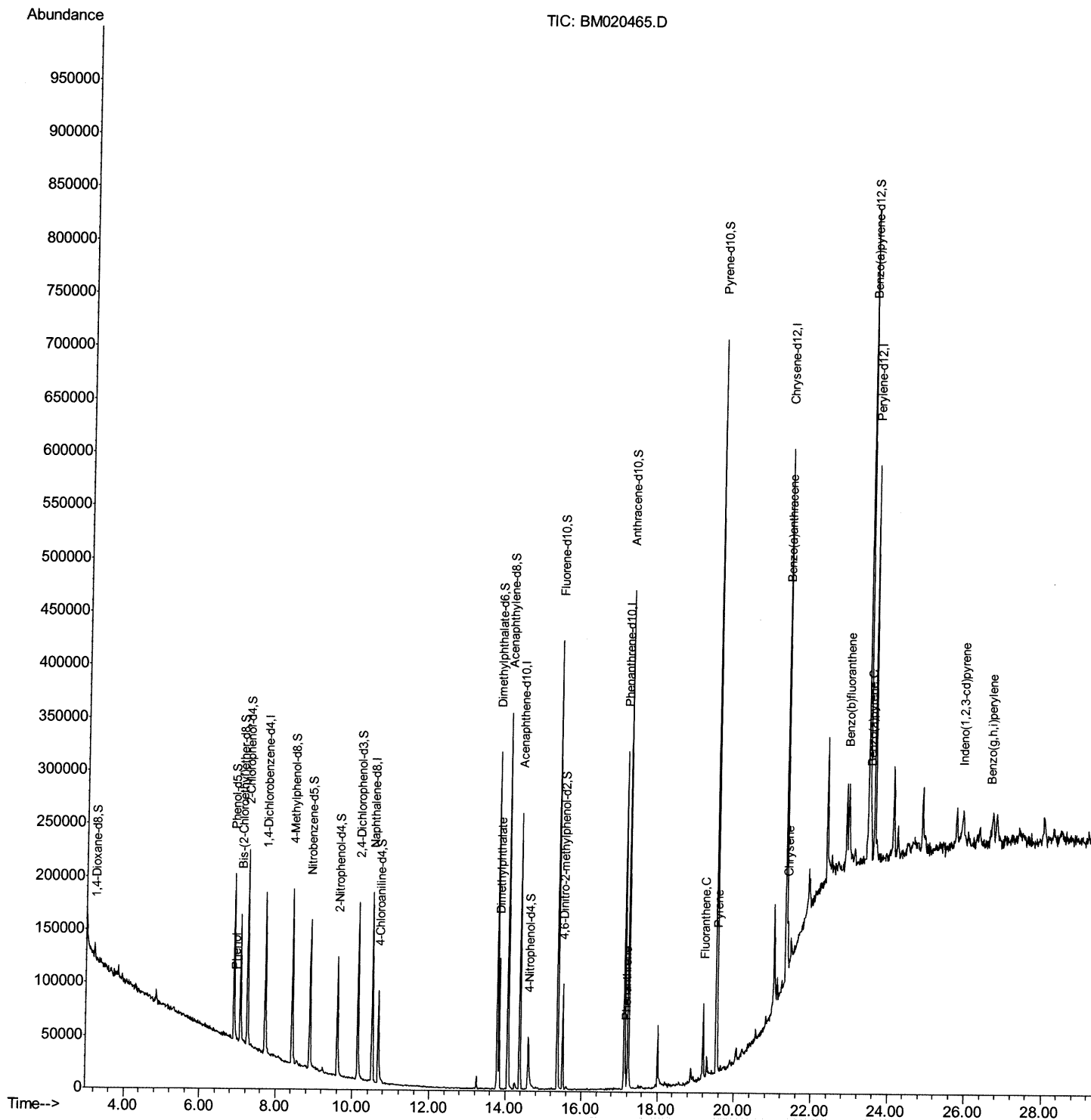
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052119\
Data File : BM020465.D
Acq On : 21 May 2019 18:18
Operator : JU/SJ
Sample : K2923-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4248

Manual Integrations
APPROVED

mohammad
5/23/2019 8:43:01 AM

Quant Time: May 22 05:31:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051619MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 22 04:54:12 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

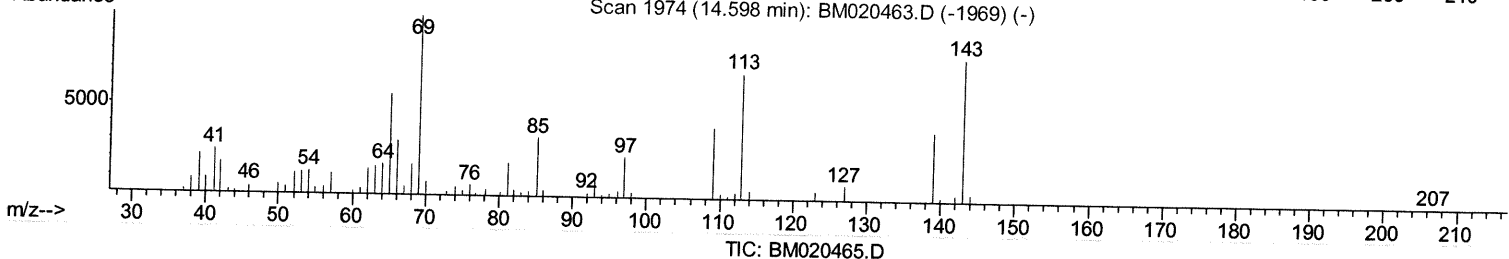
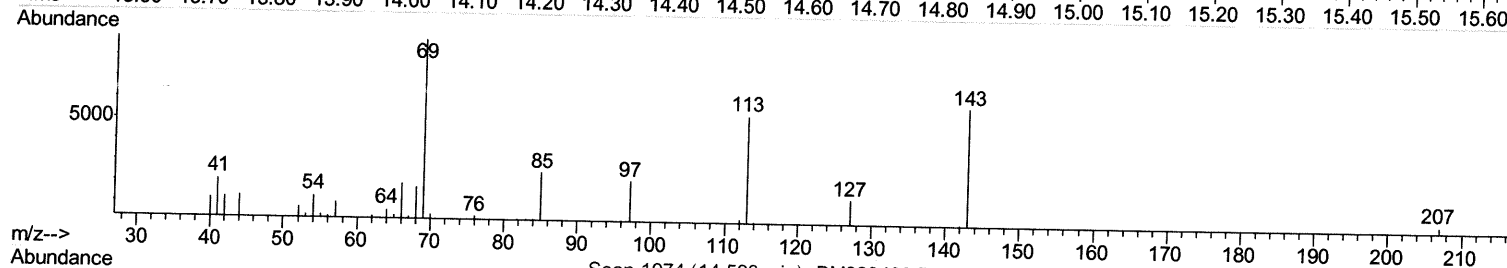
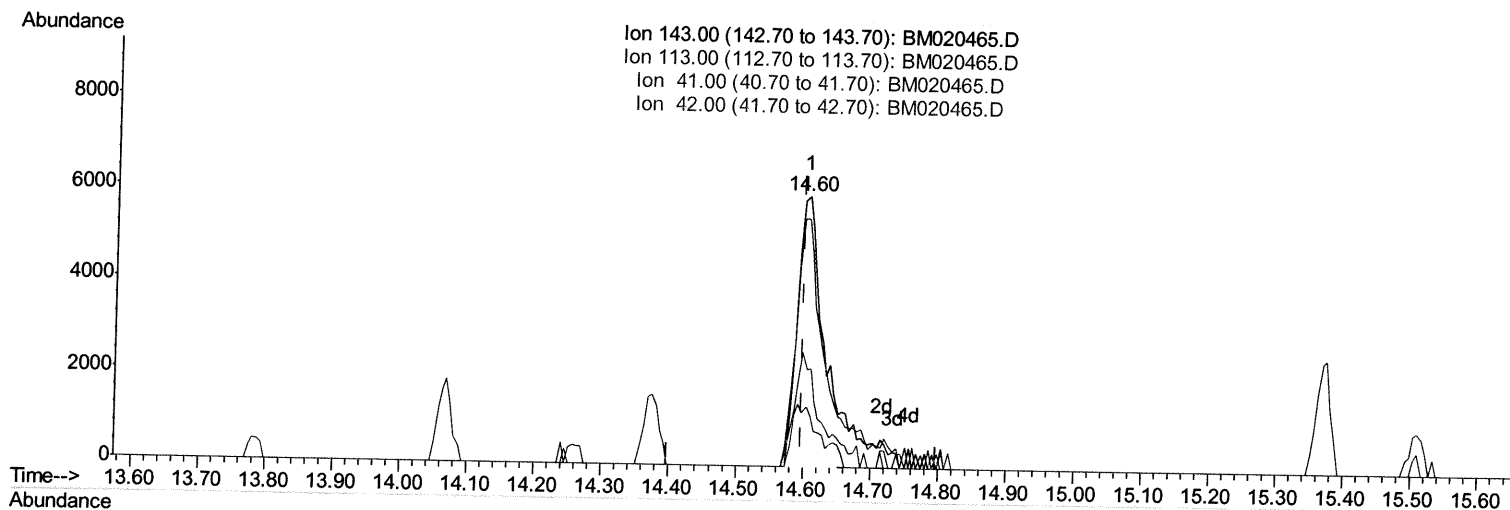
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(51) 4-Nitrophenol-d4 (S)

14.604min (+0.006) 17.42ng/ul

response 15109

Ion	Exp%	Act%
143.00	100	100
113.00	99.80	91.75
41.00	31.70	36.06
42.00	21.50	22.12

Quantitation Report (Qedit)

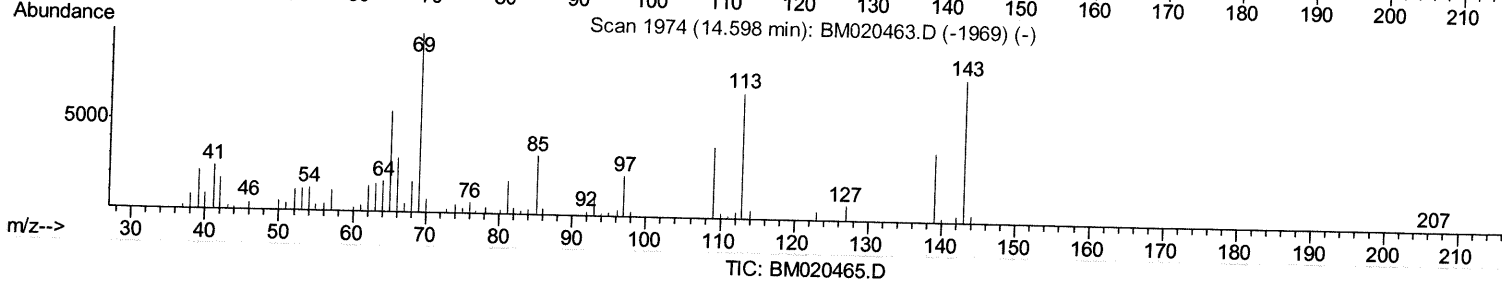
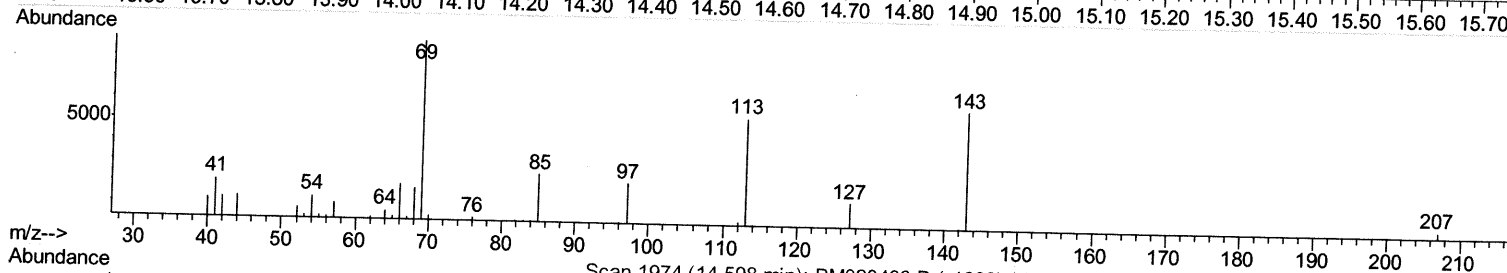
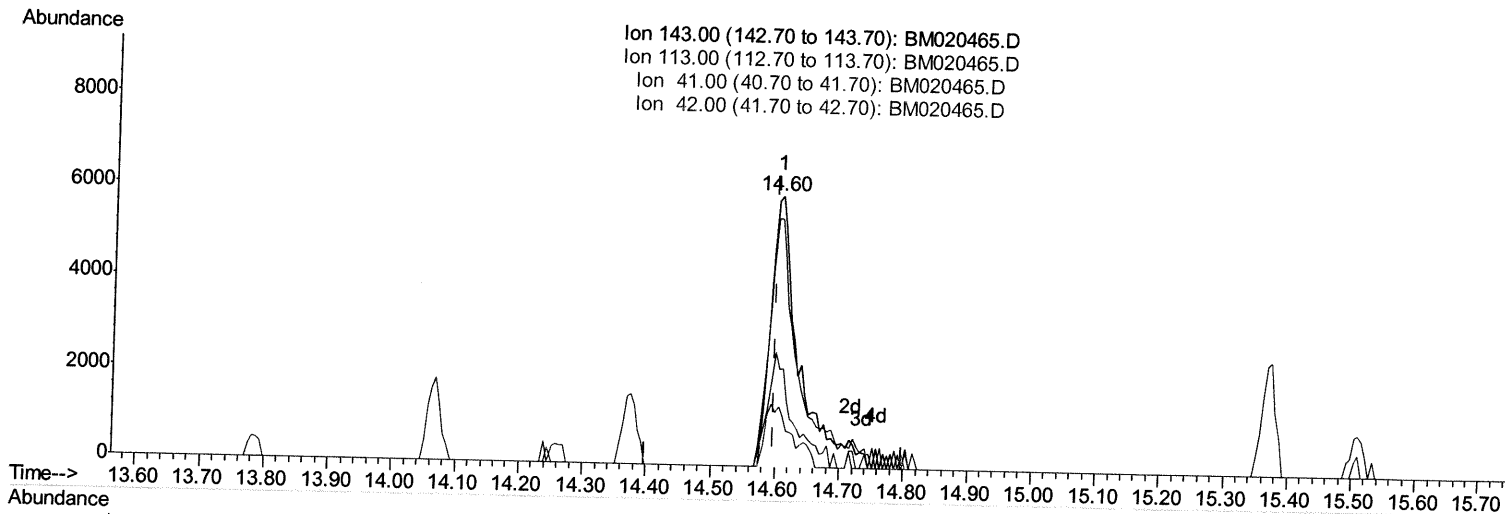
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 BNA_M
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Manual Integrations
 APPROVED

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 5/23/2019 8:43:01 AM

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(51) 4-Nitrophenol-d4 (S)

14.604min (+0.006) 21.31ng/ul m *JU* 05/24/19

response 18477

Ion	Exp%	Act%
143.00	100	100
113.00	99.80	91.75
41.00	31.70	36.06
42.00	21.50	22.12

Quantitation Report (Qedit)

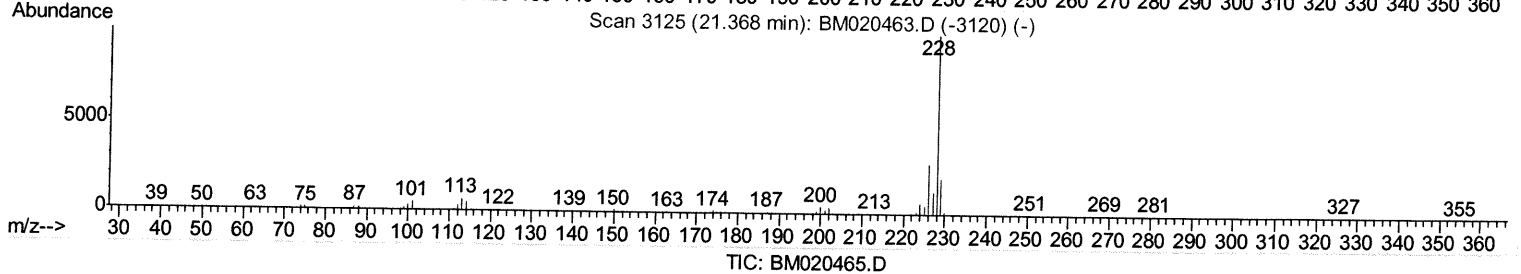
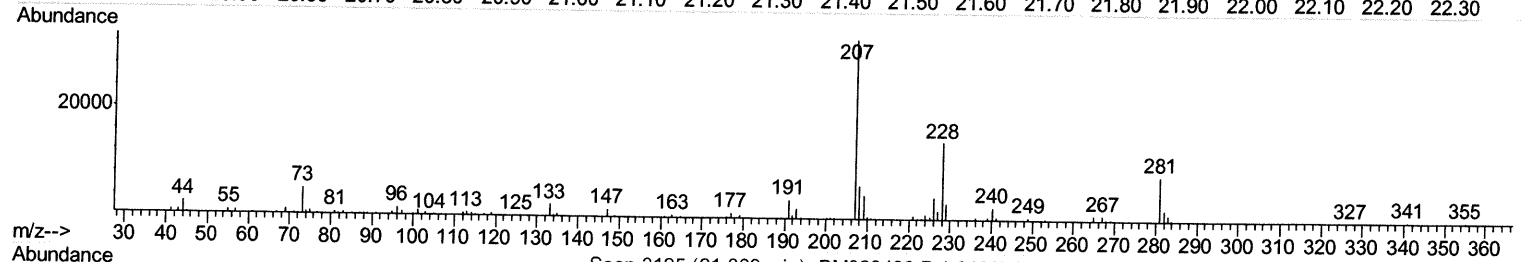
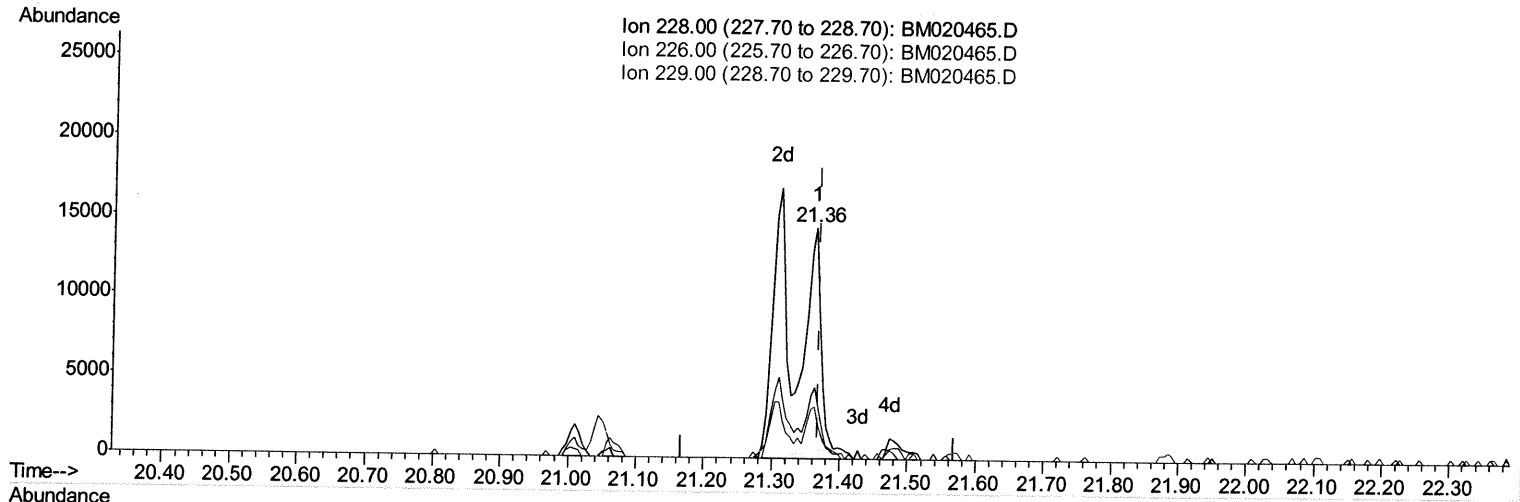
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Data File : BM020465.D
Acq On : 21 May 2019 18:18
Operator : JU/SJ
Sample : K2923-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
A4248

Manual Integrations
APPROVED

mohammad
5/23/2019 8:43:01 AM

Quant Time: May 22 04:57:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051619MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 22 04:54:12 2019
Response via : Initial Calibration



(84) Chrysene

21.362min (-0.006) 1.71ng/ul

response 22335

Ion	Exp%	Act%
228.00	100	100
226.00	27.80	30.94
229.00	19.30	22.54
0.00	0.00	0.00

Quantitation Report (Qedit)

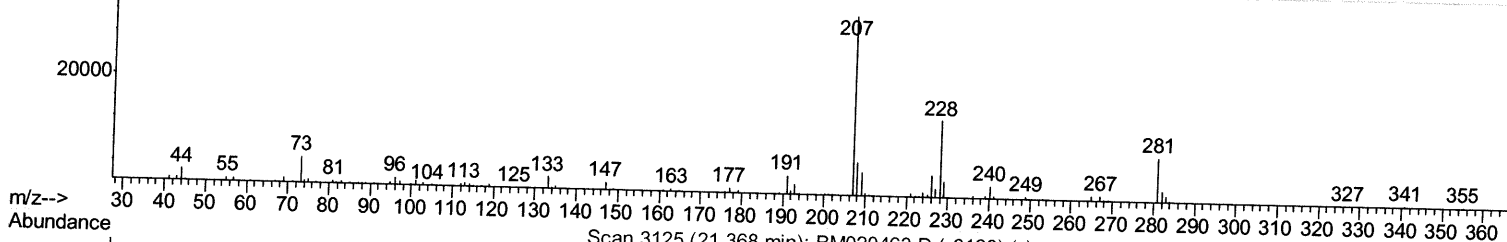
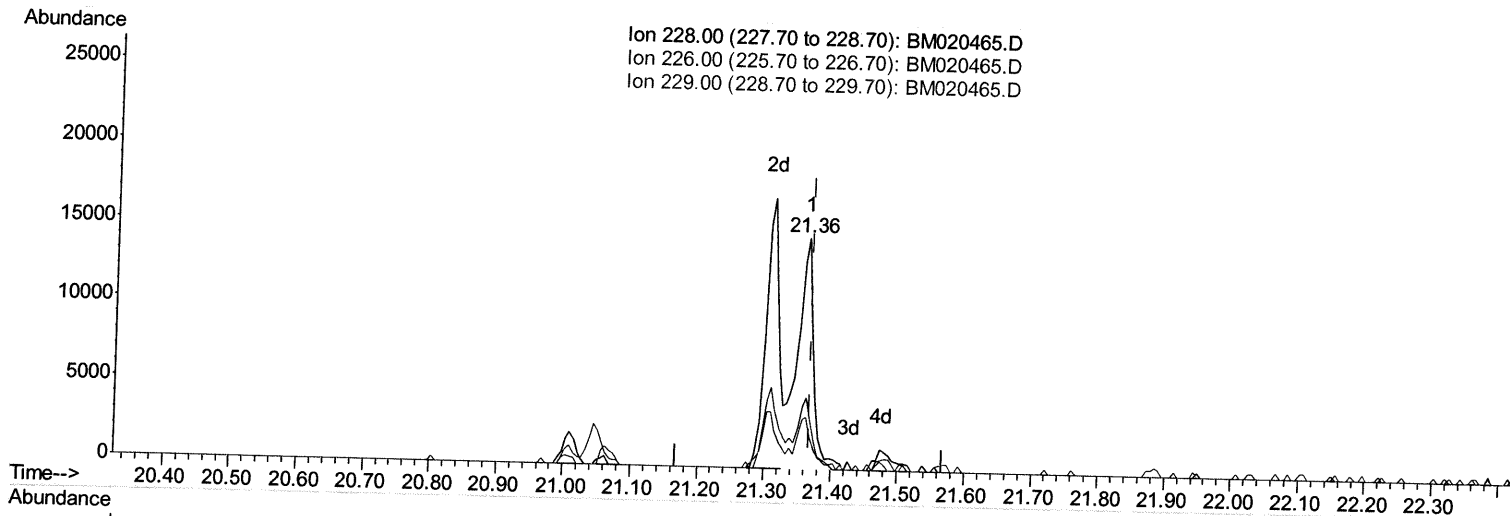
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 Data File : BM020465.D
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 Operator : JU/SJ
 Sample : K2923-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4248

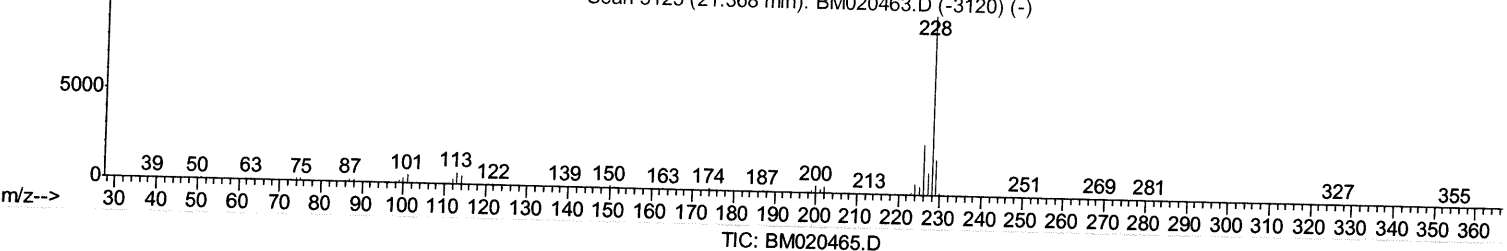
Manual Integrations
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mohammad
 5/23/2019 8:43:01 AM

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Scan 3125 (21.368 min): BM020463.D (-3120) (-)



TIC: BM020465.D

(84) Chrysene

21.362min (-0.006) 1.90ng/ul m ⁵⁰_{05/24/19}

response 24843

Ion	Exp%	Act%
228.00	100	100
226.00	27.80	30.94
229.00	19.30	22.54
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052119\
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 Acq On : 21 May 2019 18:18
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 Sample : K2923-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	42216	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	150908	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	83002	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.13	188	195570	20.00	ng/ul	0.00
77) Chrysene-d12	21.32	240	227379	20.00	ng/ul	0.00
85) Perylene-d12	23.60	264	274004	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	7795	6.81	ng/uL	0.00
5) Phenol-d5	6.89	99	101397	30.11	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	62384	29.14	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	78069	31.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	65990	26.91	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	32034	32.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	34722	32.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.14	165	69446	30.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.67	131	55438	23.44	ng/ul	0.00
43) Dimethylphthalate-d6	13.79	166	197156	33.02	ng/ul	0.00
46) Acenaphthylene-d8	14.07	160	251444	33.66	ng/ul	0.00
51) 4-Nitrophenol-d4	14.60	143	18477m	21.31	ng/ul	0.00
57) Fluorene-d10	15.37	176	176982	33.32	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	22042	19.79	ng/ul	0.00
70) Anthracene-d10	17.23	188	280713	33.46	ng/ul	0.00
78) Pyrene-d10	19.53	212	364690	33.30	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.45	264	433873	34.84	ng/ul	0.00

Target Compounds

					Qvalue	
6) Phenol	6.92	94	3923	1.107	ng/ul#	87
44) Dimethylphthalate	13.83	163	73555	12.451	ng/ul	98
69) Phenanthrene	17.17	178	14174	1.521	ng/ul	97
76) Fluoranthene	19.19	202	41952	3.564	ng/ul	99
79) Pyrene	19.56	202	40650	2.979	ng/ul	99
82) Benzo(a)anthracene	21.31	228	23228	1.697	ng/ul	94
84) Chrysene	21.36	228	24843m	1.900	ng/ul	94
87) Benzo(b)fluoranthene	22.91	252	37346	2.388	ng/ul#	91
90) Benzo(a)pyrene	23.50	252	24300	1.651	ng/ul#	95
91) Indeno(1,2,3-cd)pyrene	25.90	276	21702	1.294	ng/ul	96
93) Benzo(g,h,i)perylene	26.61	276	17325	1.248	ng/ul#	93

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