

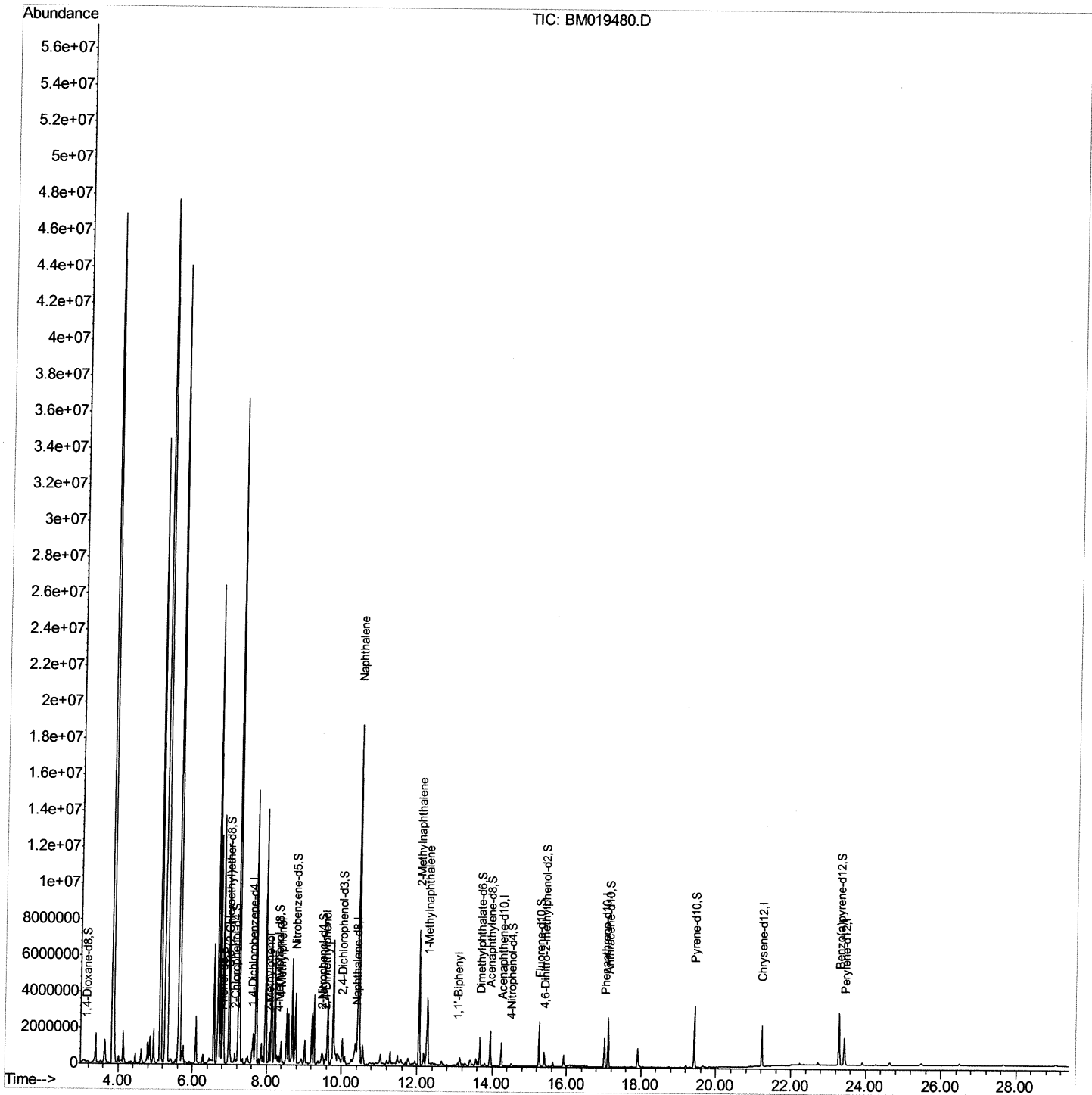
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019480.D
Acq On : 26 Mar 2019 21:05
Operator : JU/SJ
Sample : K2063-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB6R9

Manual Integrations
APPROVED

Sohil
3/27/2019 5:46:32 PM

Quant Time: Mar 27 04:16:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 27 03:30:35 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

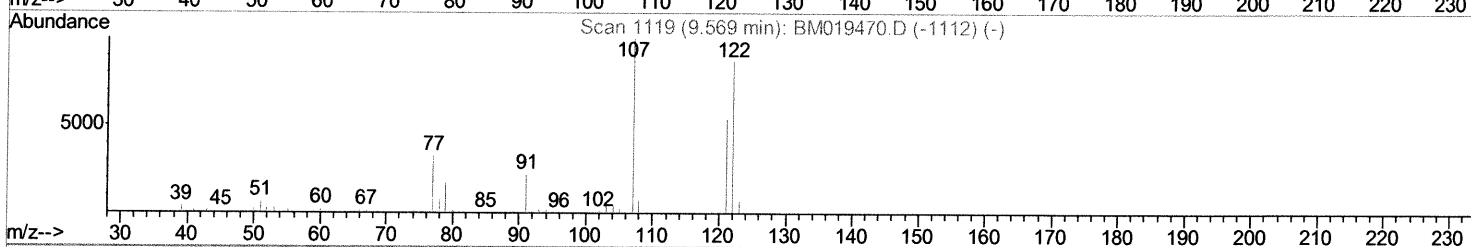
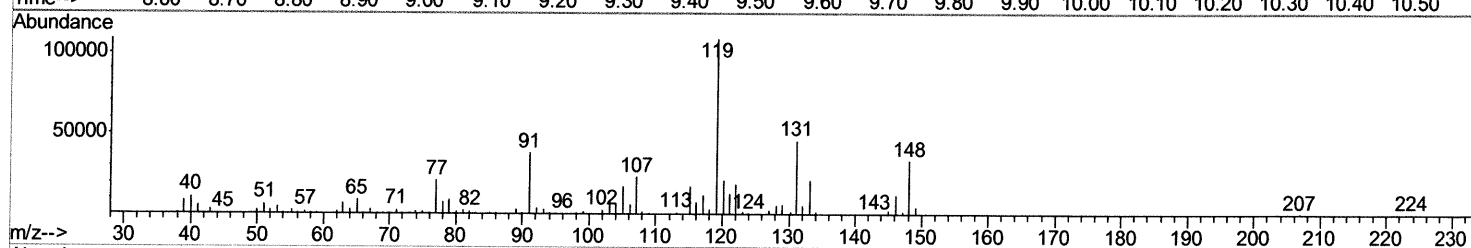
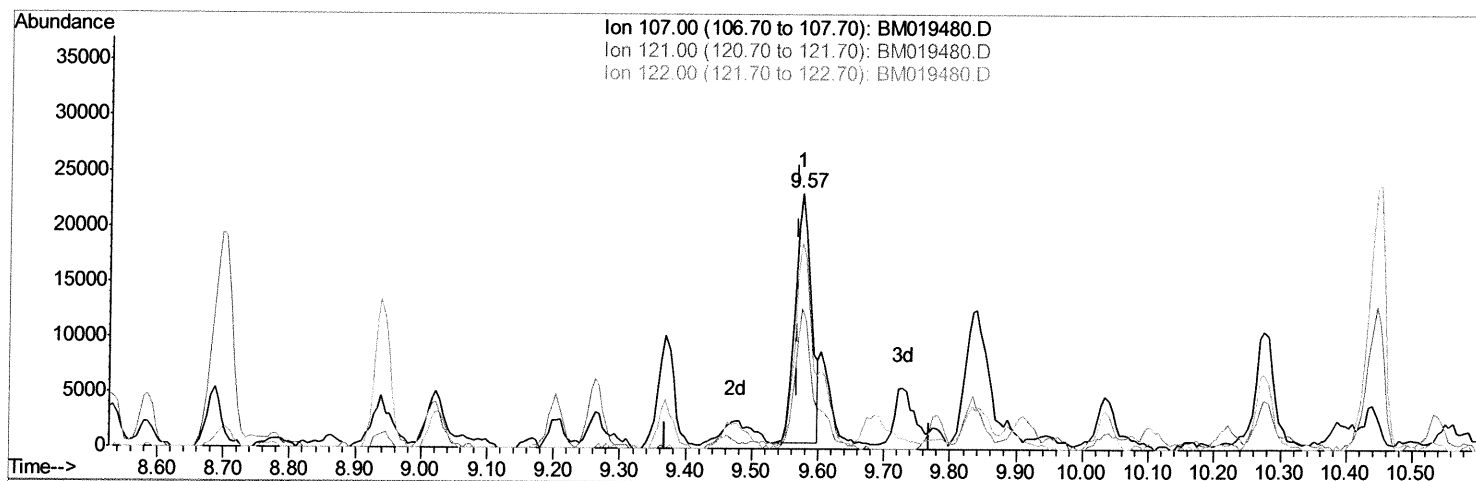
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TIC: BM019480.D

(24) 2,4-Dimethylphenol

9.575min (+0.006) 2.92ng/ul

response 38527

Ion	Exp%	Act%
107.00	100	100
121.00	47.60	54.69
122.00	79.80	80.49
0.00	0.00	0.00

Quantitation Report (Qedit)

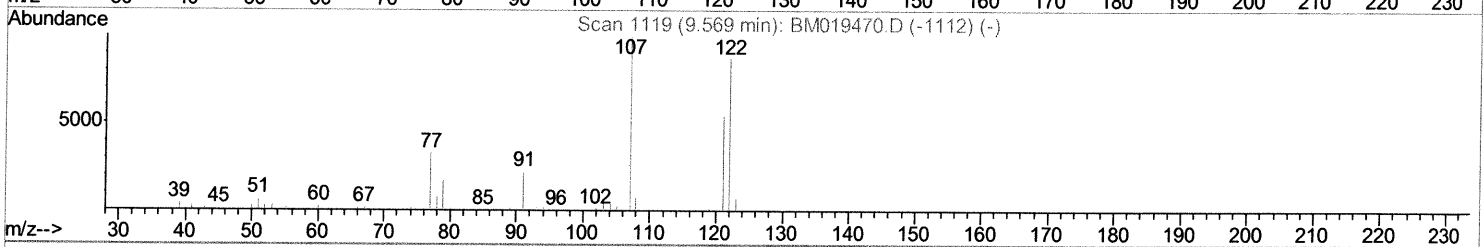
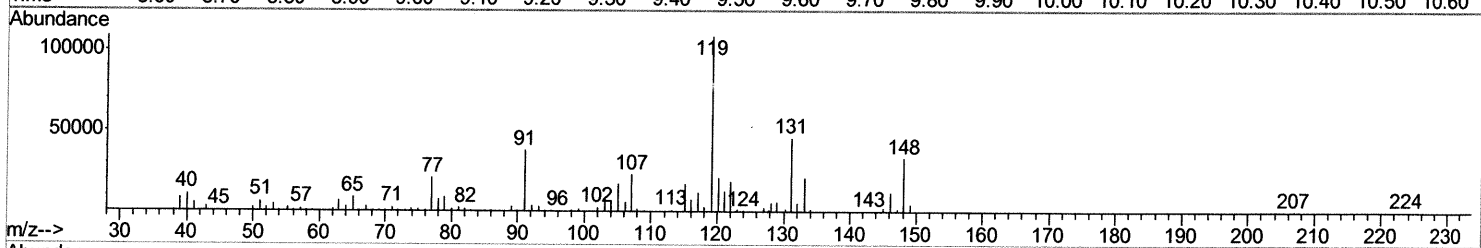
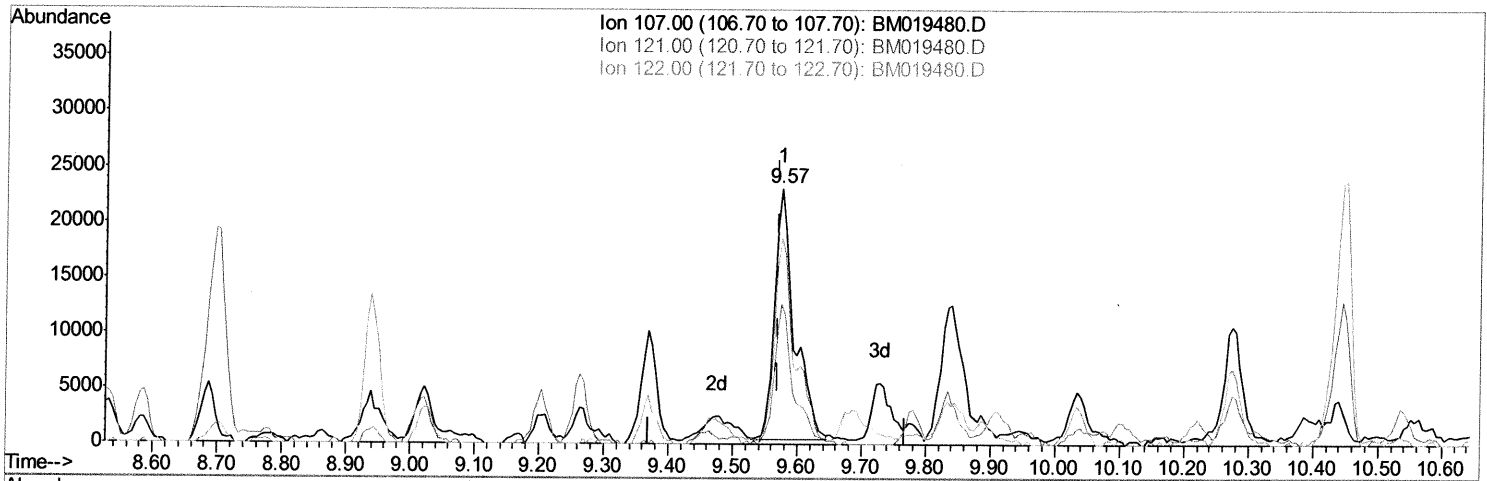
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Manual Integrations
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TIC: BM019480.D

(24) 2,4-Dimethylphenol

9.575min (+0.006) 3.69ng/ul m SD 4/2/19

response 48659

Ion	Exp%	Act%
107.00	100	100
121.00	47.60	54.69
122.00	79.80	80.49
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
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 Sample : K2063-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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 BNA_M
 ClientSampled :
 DB6R9

Manual Integrations
 APPROVED

Sohil
 3/27/2019 5:46:32 PM

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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 27 03:30:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	166565	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	693542	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	400651	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	904436	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	1017615	20.00	ng/ul	0.00
86) Perylene-d12	23.43	264	1043814	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.15	96	12073	2.84	ng/uL	0.00
5) Phenol-d5	6.79	99	106104	7.15	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	265928	26.78	ng/ul	0.01
9) 2-Chlorophenol-d4	7.14	132	290235	25.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	194593	17.48	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	227388	45.96	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	193330	35.11	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	337857	32.52	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	7862	0.63	ng/ul	0.04
44) Dimethylphthalate-d6	13.67	166	1078445	33.36	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	1339180	33.13	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	39097	7.72	ng/ul	0.02
58) Fluorene-d10	15.26	176	935040	33.57	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	189110	31.62	ng/ul	0.00
71) Anthracene-d10	17.12	188	1542733	35.55	ng/ul	0.00
79) Pyrene-d10	19.42	212	1728172	36.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.29	264	1841286	33.27	ng/ul	0.00
Target Compounds						
11) 2-Methylphenol	8.05	108	26088	2.417	ng/ul	97
16) 4-Methylphenol	8.38	108	33884	2.862	ng/ul	89
24) 2,4-Dimethylphenol	9.57	107	48659m	3.687	ng/ul	
28) Naphthalene	10.45	128	15707474	437.188	ng/ul	96
34) 2-Methylnaphthalene	12.06	142	3511004	139.463	ng/ul	100
35) 1-Methylnaphthalene	12.28	142	1631877	68.651	ng/ul	96
41) 1,1'-Biphenyl	13.08	154	62349	1.827	ng/ul#	91

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

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 4/2/19