

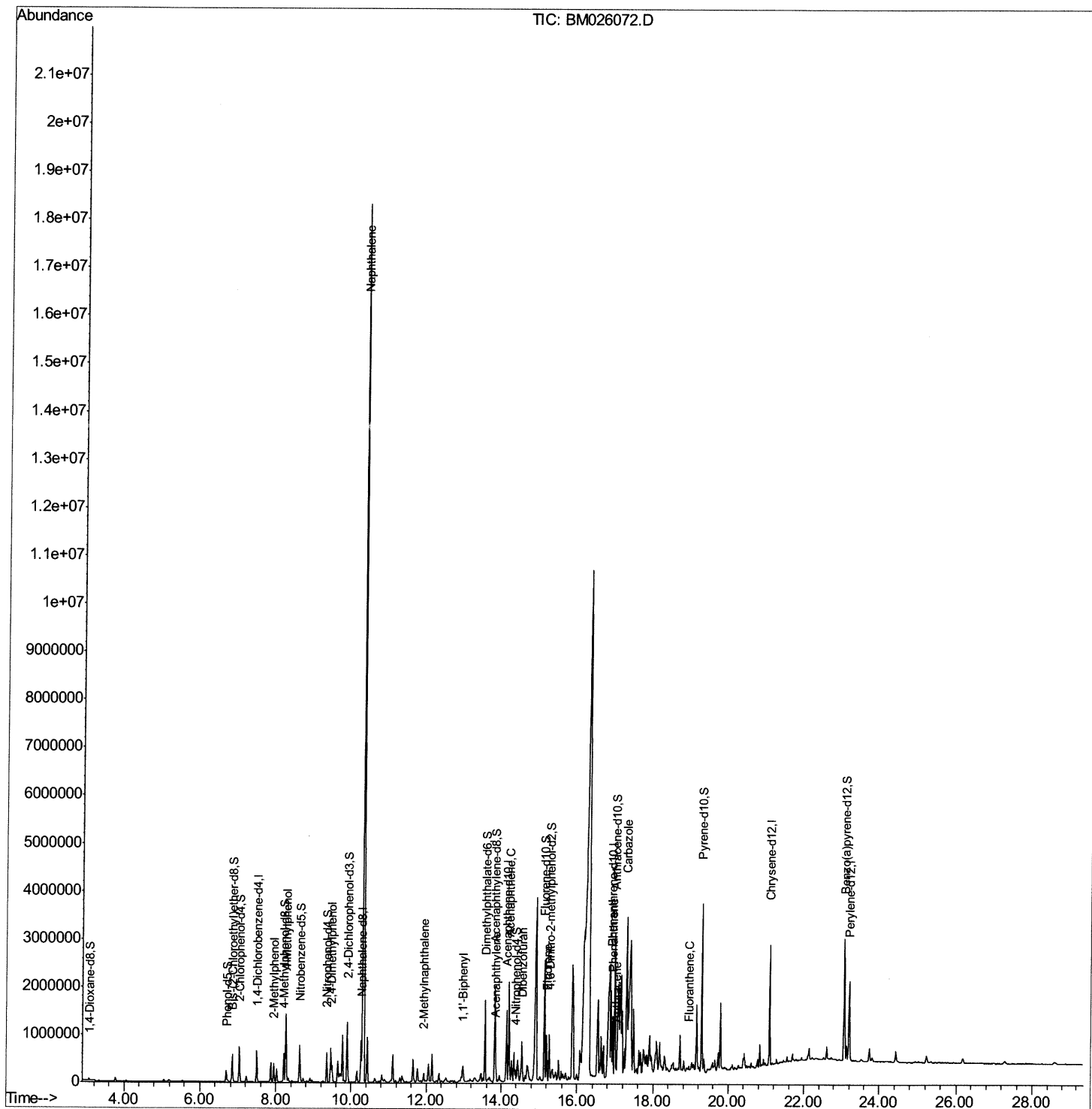
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052120\
Data File : BM026072.D
Acq On : 22 May 2020 00:22
Operator : MA/SJ
Sample : L2725-12
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9B23

Manual Integrations
APPROVED

mohammad
5/22/2020 2:54:54 PM

Quant Time: May 22 05:29:07 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 21 10:18:34 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

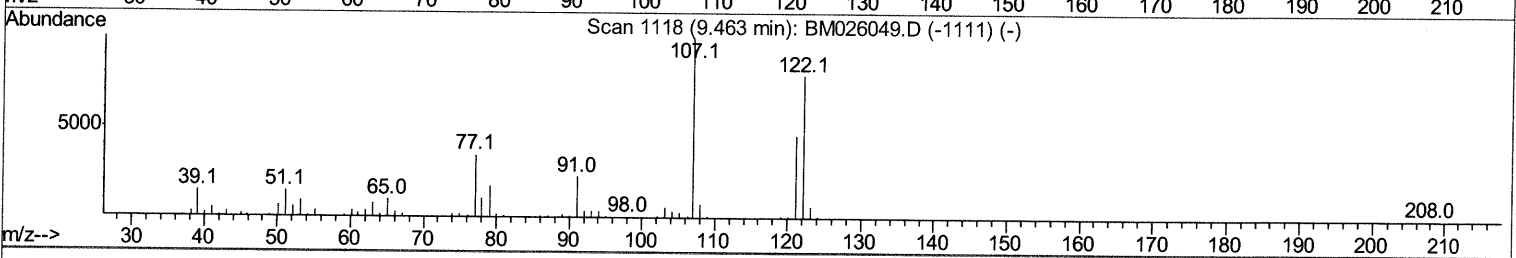
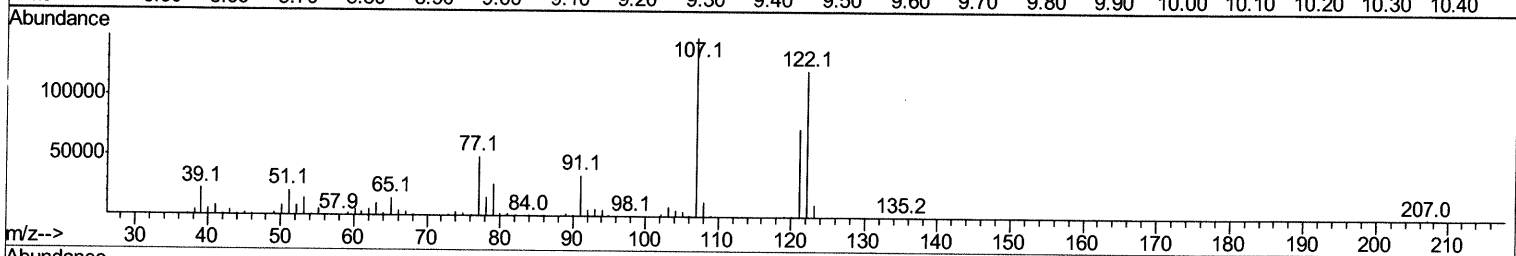
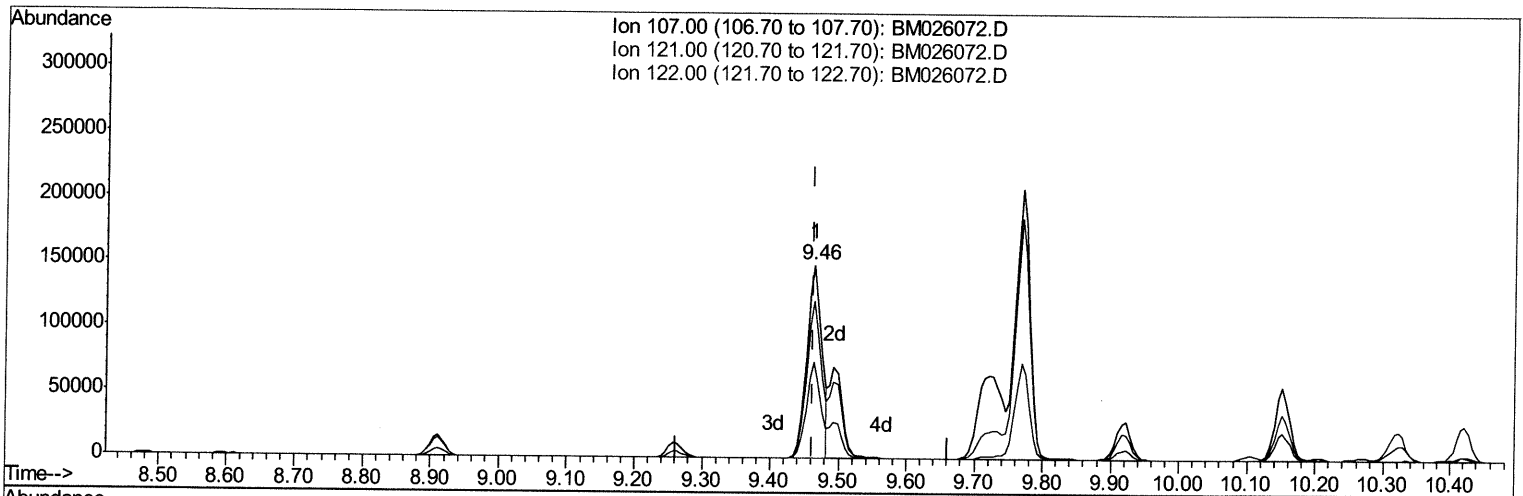
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TIC: BM026072.D

(24) 2,4-Dimethylphenol

9.463min (+0.000) 14.70ng/ul

response 227090

Ion	Exp%	Act%
107.00	100	100
121.00	46.90	49.39
122.00	79.70	81.63
0.00	0.00	0.00

Quantitation Report (Qedit)

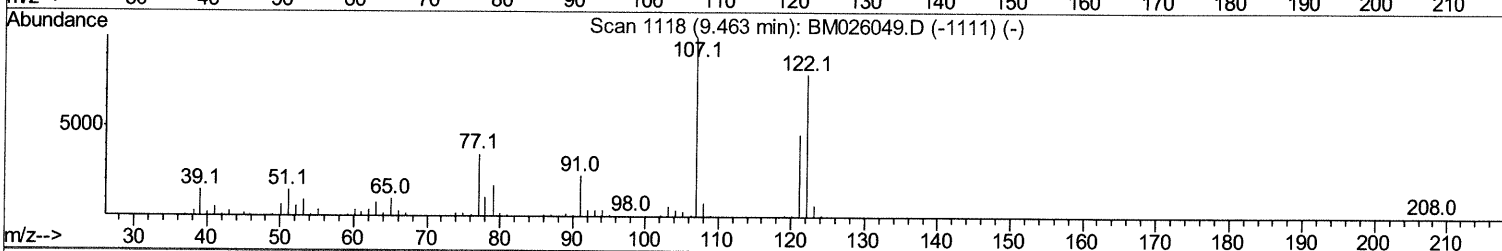
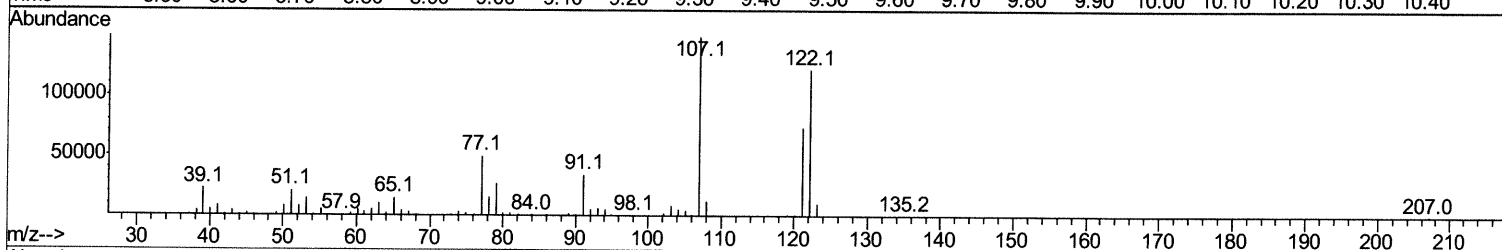
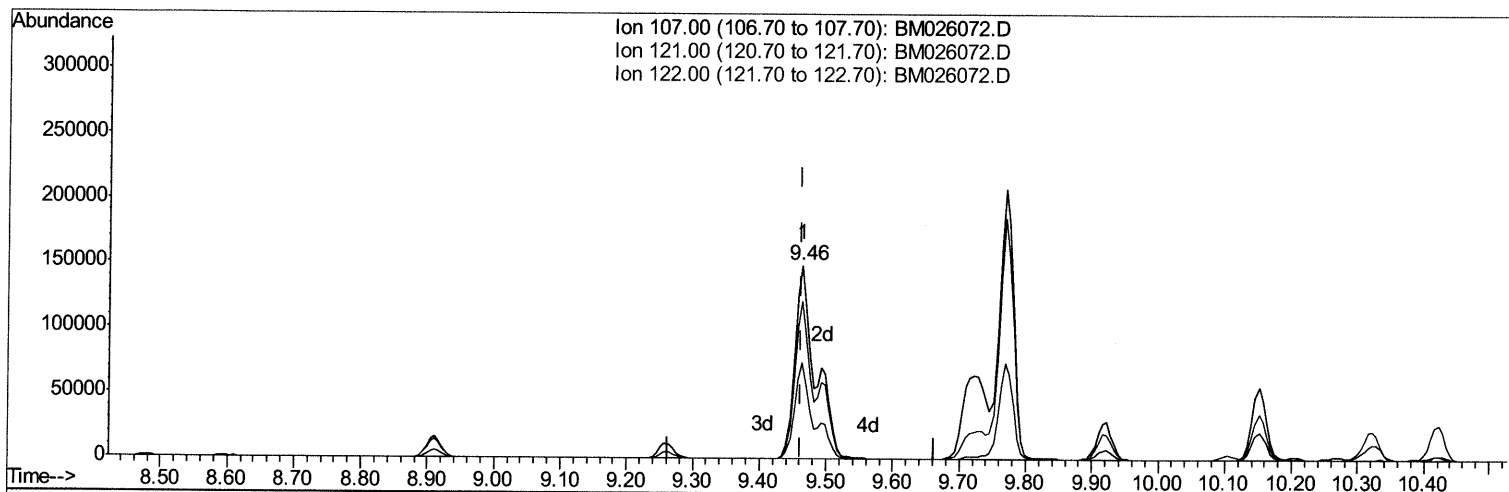
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(24) 2,4-Dimethylphenol

9.463min (+0.000) 20.81ng/ul m

05/22/20 AP

response 321433

Ion	Exp%	Act%
107.00	100	100
121.00	46.90	49.39
122.00	79.70	81.63
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052120\
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Manual Integrations
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mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	158526	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	689697	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	431871	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	957377	20.00	ng/ul	0.00
78) Chrysene-d12	21.09	240	1031104	20.00	ng/ul	0.00
86) Perylene-d12	23.21	264	1018062	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.07	96	21421	4.08	ng/uL	0.00
5) Phenol-d5	6.69	99	112831	7.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	260933	26.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.03	132	277008	24.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.21	113	204315	18.85	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	152098	27.05	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	167737	30.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	330164	29.57	ng/ul	0.00
29) 4-Chloroaniline-d4	10.41	131	7958	0.70	ng/ul	0.01
44) Dimethylphthalate-d6	13.56	166	1025051	29.60	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	1211622	29.53	ng/ul	0.00
52) 4-Nitrophenol-d4	14.38	143	48783	7.95	ng/ul	0.02
58) Fluorene-d10	15.14	176	898589	29.83	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.27	200	163885	29.36	ng/ul	0.00
71) Anthracene-d10	16.99	188	1433511	30.39	ng/ul	0.00
79) Pyrene-d10	19.29	212	1642582	31.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.07	264	1511869	27.90	ng/ul	0.00
Target Compounds						
					Ovalue	
11) 2-Methylphenol	7.95	108	114194	10.281	ng/ul	100
16) 4-Methylphenol	8.27	108	493383	40.844	ng/ul	89
24) 2,4-Dimethylphenol	9.46	107	321433m	20.811	ng/ul	89
28) Naphthalene	10.32	128	14526426	344.478	ng/ul	99
34) 2-Methylnaphthalene	11.93	142	80022	2.821	ng/ul	92
41) 1,1'-Biphenyl	12.96	154	135438	3.534	ng/ul	98
48) Acenaphthylene	13.86	152	49671	1.103	ng/ul#	89
50) Acenaphthene	14.20	153	740037	23.295	ng/ul	98
54) Dibenzofuran	14.55	168	455453	10.202	ng/ul	99
59) Fluorene	15.20	166	372443	10.313	ng/ul	100
70) Phenanthrene	16.93	178	617612	10.241	ng/ul	100
72) Anthracene	17.02	178	75093	1.241	ng/ul	98
75) Carbazole	17.30	167	1486256	29.693	ng/ul	100
77) Fluoranthene	18.96	202	63748	1.008	ng/ul#	95

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