

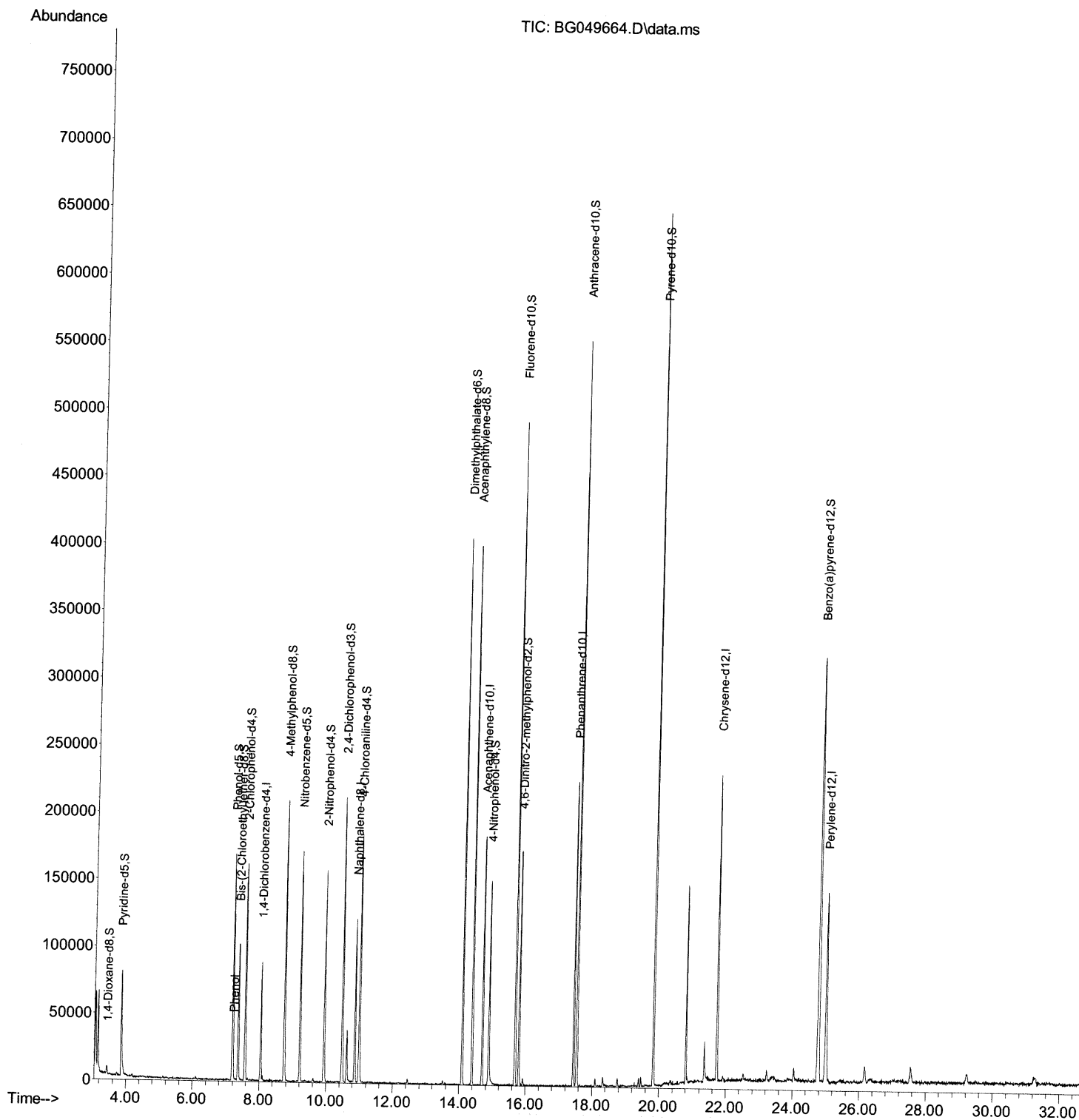
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049664.D
Acq On : 6 Aug 2021 00:51
Operator : CG/JU
Sample : M3162-12
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGEW3

Manual Integrations
APPROVED

mohammad
8/6/2021 2:48:34 PM

Quant Time: Aug 06 04:03:04 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 04 16:17:29 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

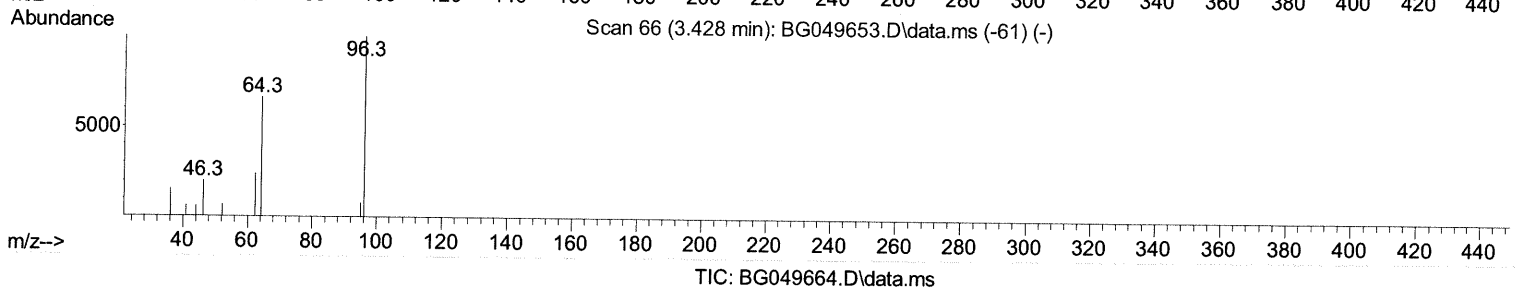
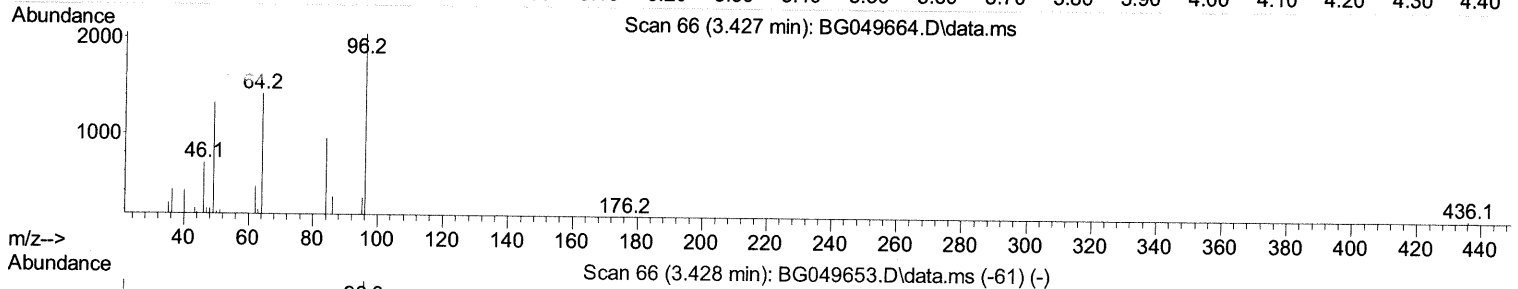
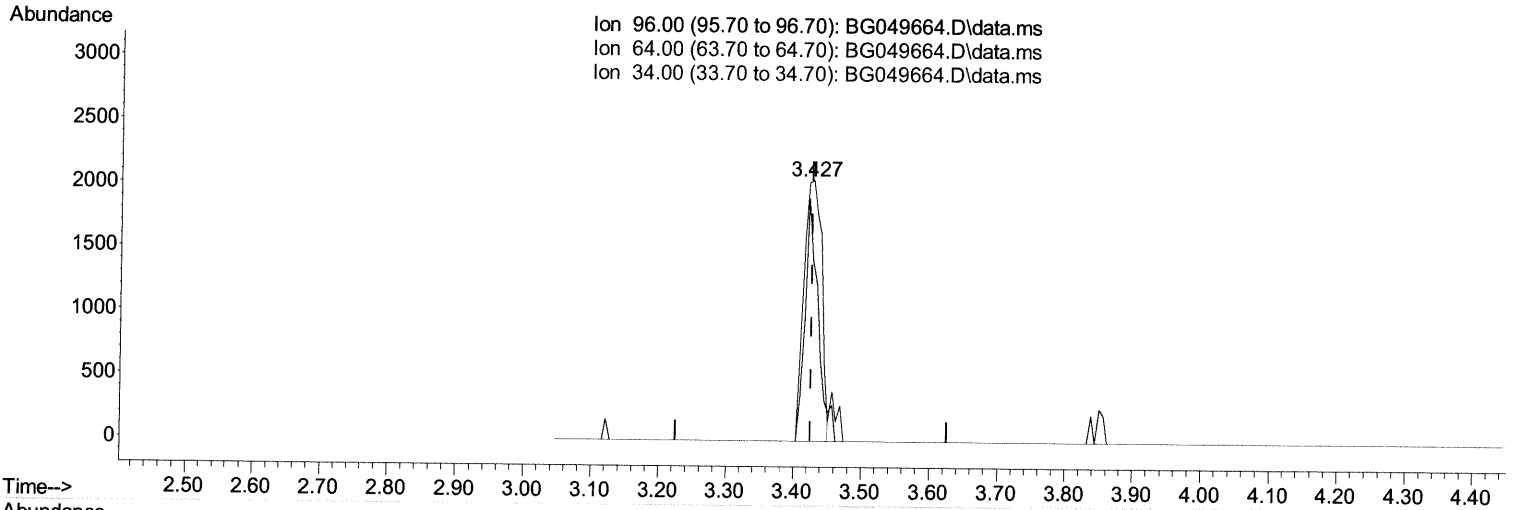
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TIC: BG049664.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.427min (+ 0.001) 7.10 ng/uL

response 3706

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.10	69.31
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

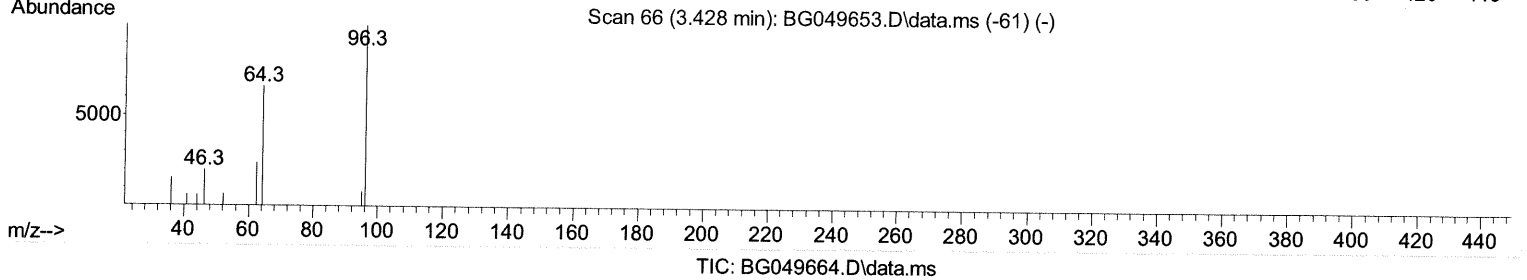
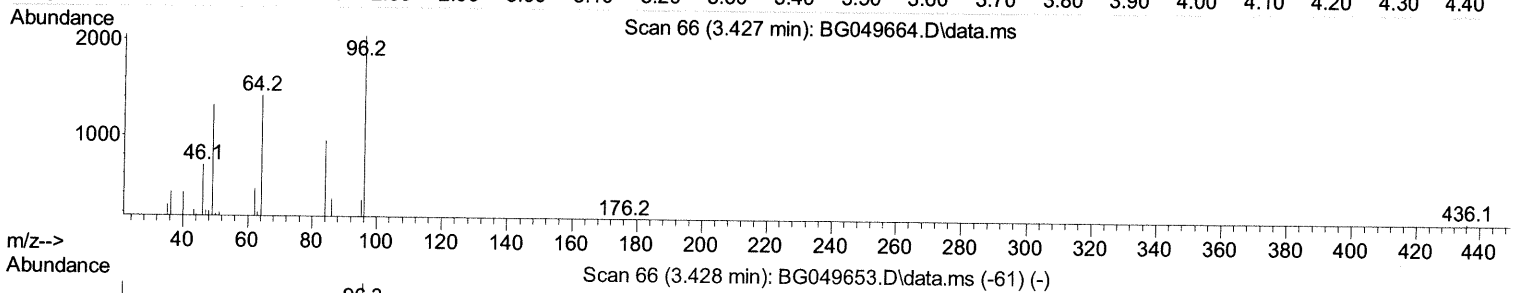
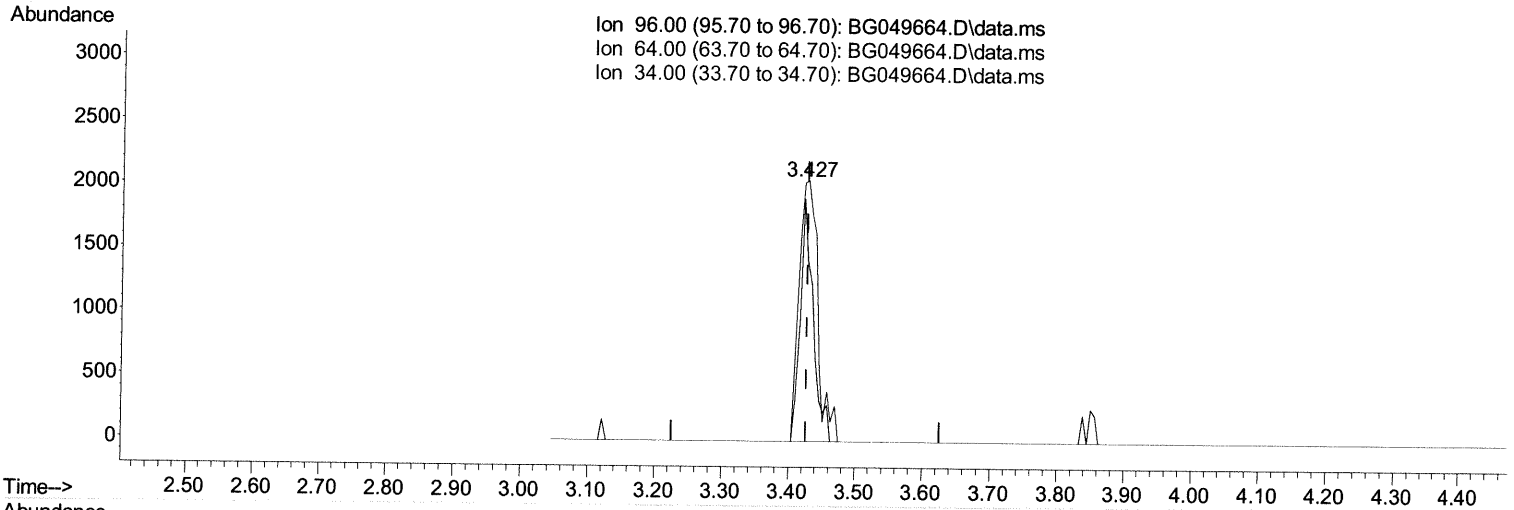
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(3) 1,4-Dioxane-d8 (S)

3.427min (+ 0.001) 7.64 ng/uL m 08/09/21 JU

response 3992

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.10	69.31
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

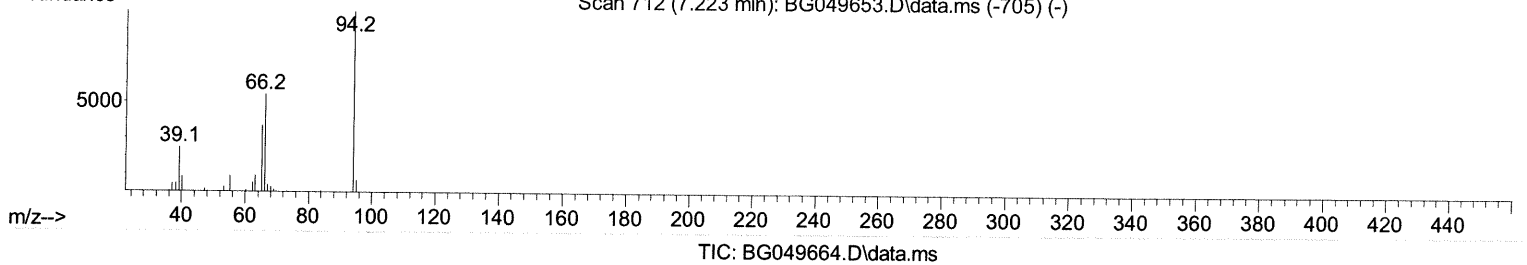
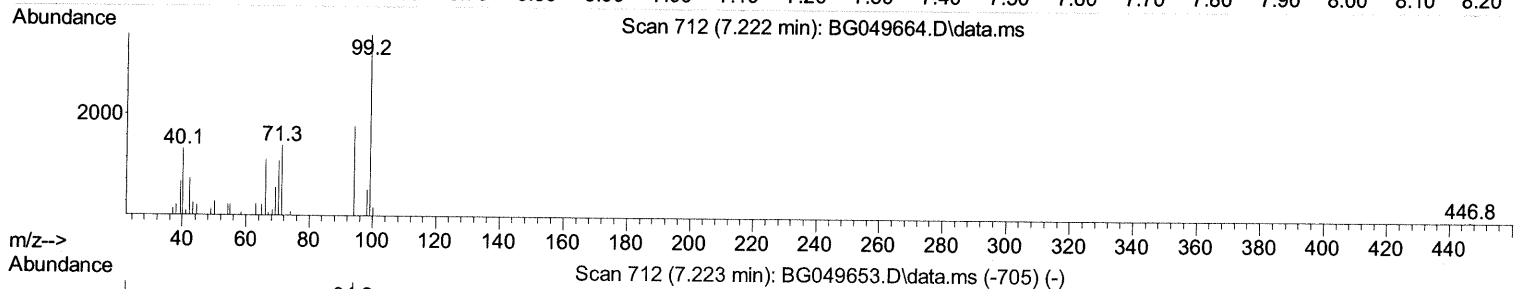
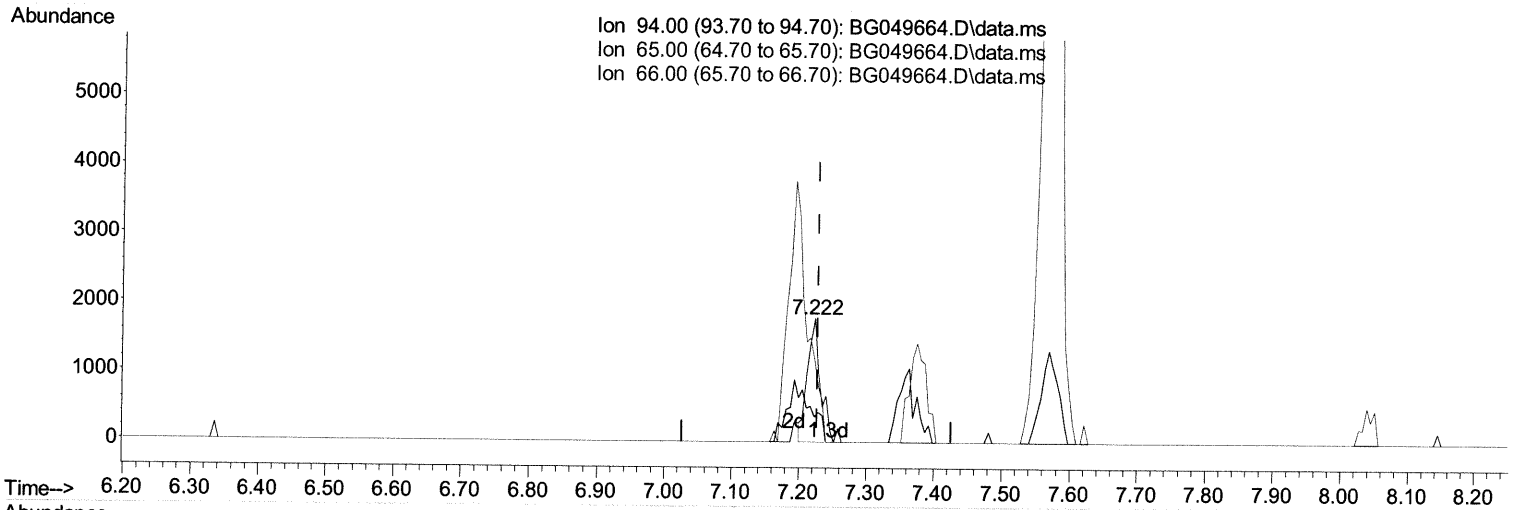
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TIC: BG049664.D\data.ms

(8) Phenol

7.222min (-0.004) 1.10 ng/ul

response 2407

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	30.90	20.28#
66.00	54.60	65.49
0.00	0.00	0.00

Quantitation Report (Qedit)

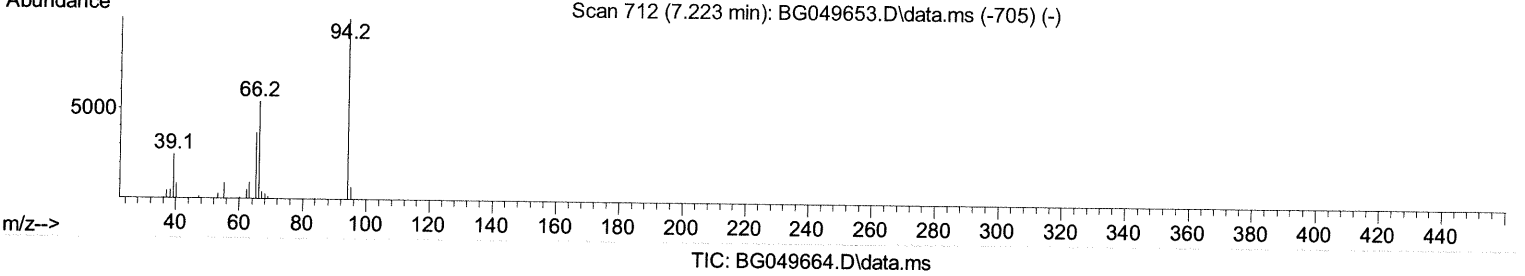
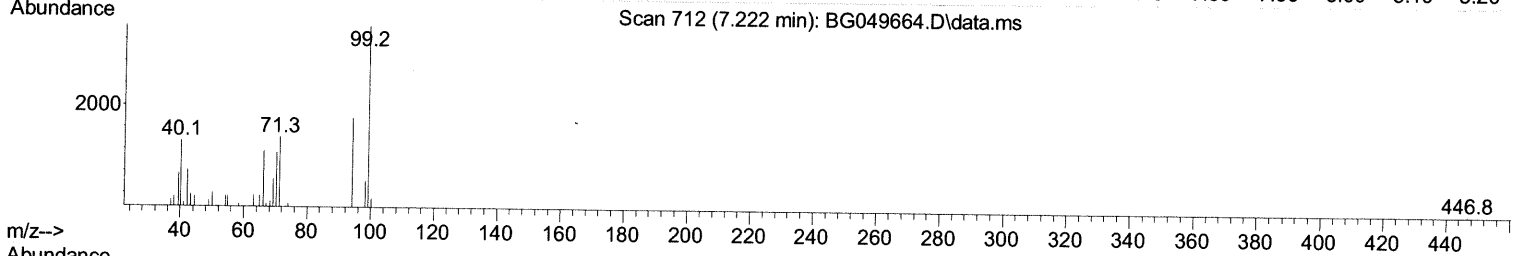
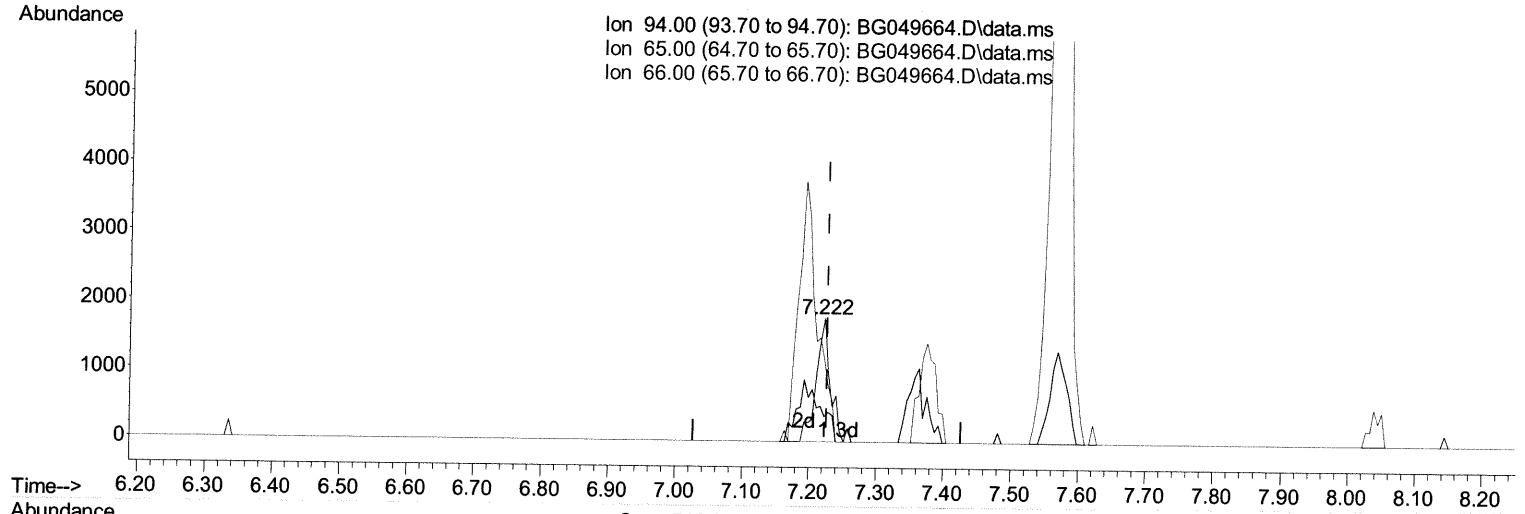
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 Sample : M3162-12
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
 APPROVED

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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



(8) Phenol

7.222min (-0.004) 1.18 ng/ul m 08/09/21 JU

response 2598

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	30.90	20.28#
66.00	54.60	65.49
0.00	0.00	0.00

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 Misc :
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Instrument :
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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.039	152	24420	20.000	ng/ul	0.00
20) Naphthalene-d8	10.864	136	98817	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.694	164	67696	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.444	188	141177	20.000	ng/ul	0.00
79) Chrysene-d12	21.720	240	139585	20.000	ng/ul	# 0.00
88) Perylene-d12	24.998	264	142927	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.427	96	3992m	7.643	ng/ul	0.00
4) Pyridine-d5	3.844	84	53038	33.839	ng/ul	0.00
7) Phenol-d5	7.193	99	90077	40.644	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	51002	40.063	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	66245	41.763	ng/ul	0.00
15) 4-Methylphenol-d8	8.744	113	72027	42.624	ng/ul	0.00
21) Nitrobenzene-d5	9.208	128	35739	45.913	ng/ul	0.00
24) 2-Nitrophenol-d4	9.936	143	40582	45.651	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.477	165	69589	41.516	ng/ul	0.00
31) 4-Chloroaniline-d4	10.999	131	95952	41.815	ng/ul	0.00
46) Dimethylphthalate-d6	14.089	166	242575	45.661	ng/ul	0.00
49) Acenaphthylene-d8	14.383	160	286281	45.295	ng/ul	0.00
54) 4-Nitrophenol-d4	14.888	143	31883	37.119	ng/ul	0.00
60) Fluorene-d10	15.687	176	199059	43.591	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.805	200	37036	38.978	ng/ul	0.00
73) Anthracene-d10	17.543	188	316802	47.539	ng/ul	0.00
81) Pyrene-d10	19.829	212	361130	46.956	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.769	264	341564	43.614	ng/ul	-0.01
Target Compounds						
8) Phenol	7.222	94	2598m	1.183	ng/ul	Qvalue

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