

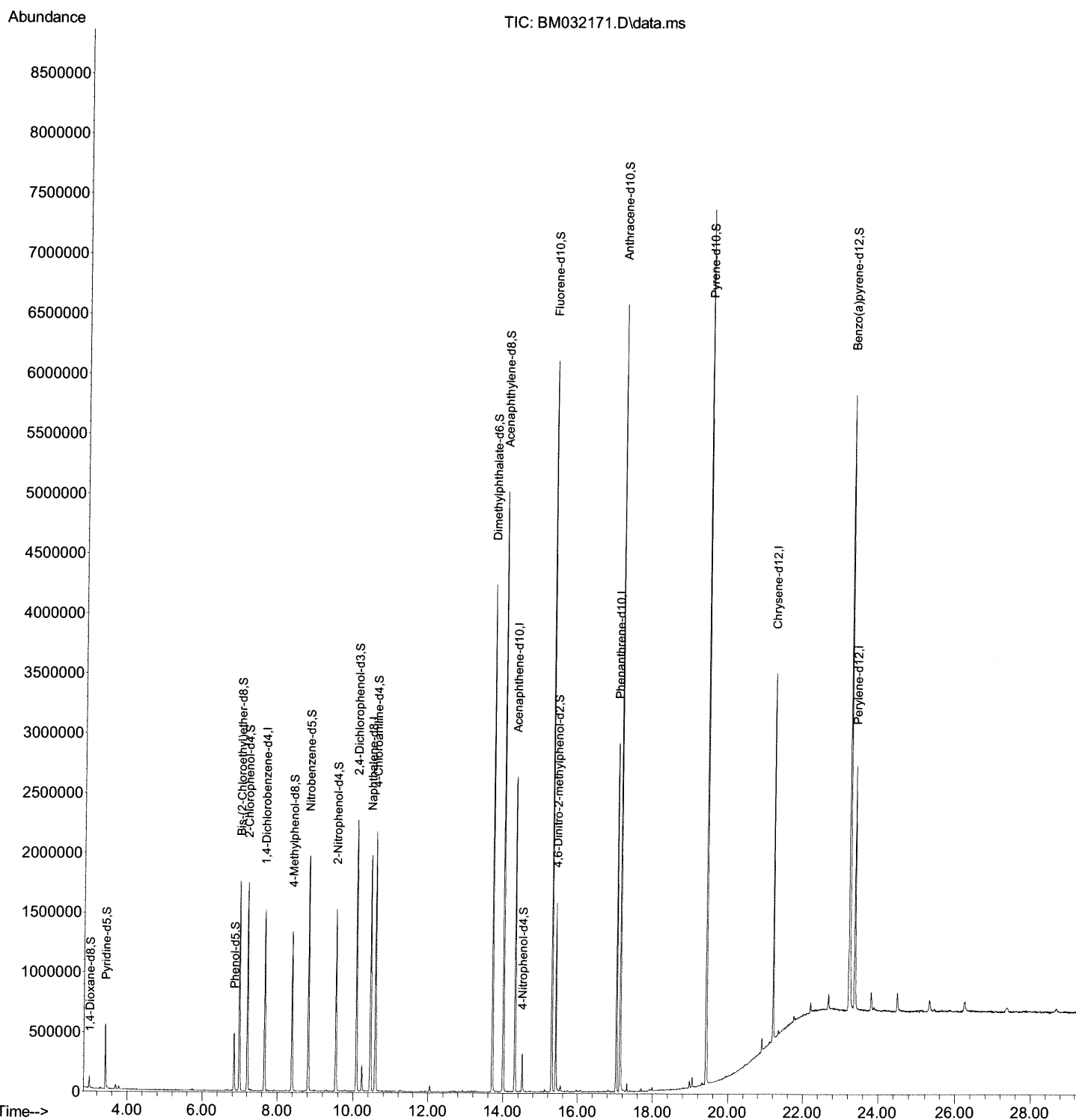
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
Data File : BM032171.D
Acq On : 26 Sep 2021 02:04
Operator : CG/JU
Sample : M3739-19
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG1

Manual Integrations
APPROVED

mohammad
9/27/2021 3:06:35 PM

Quant Time: Sep 26 03:18:55 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM092421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 25 01:15:33 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

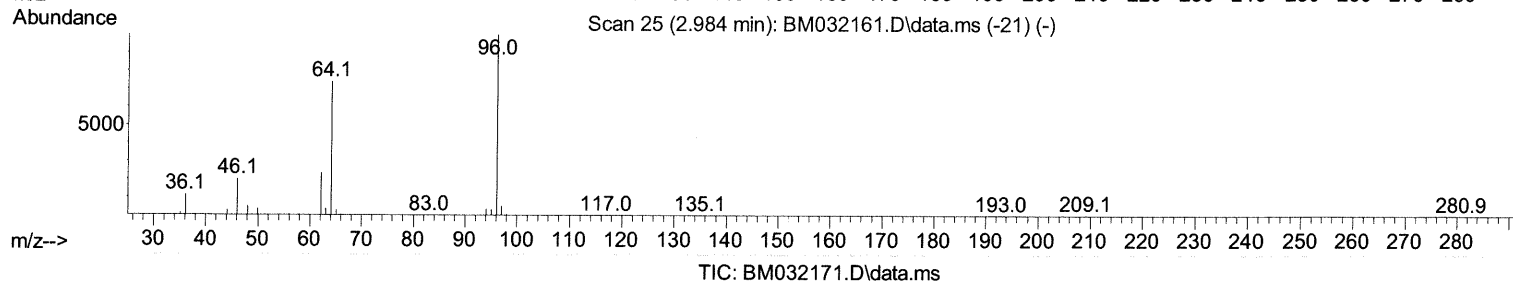
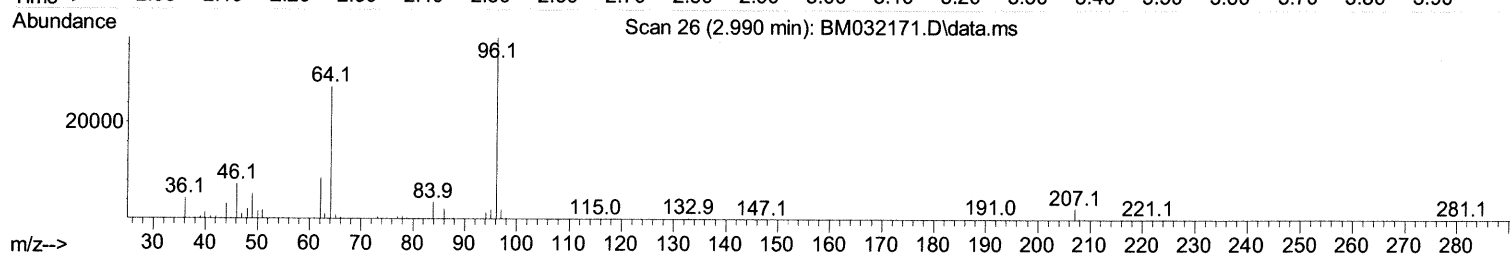
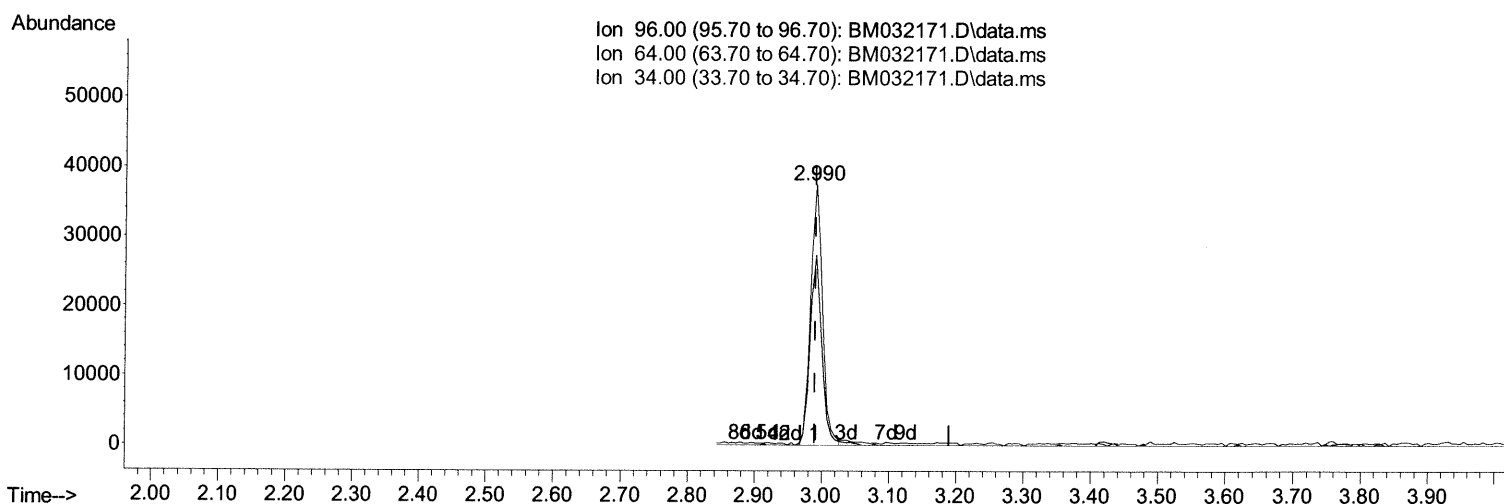
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(3) 1,4-Dioxane-d8 (S)

2.990min (-0.000) 4.83 ng/uL

response 45508

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.40	73.02
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

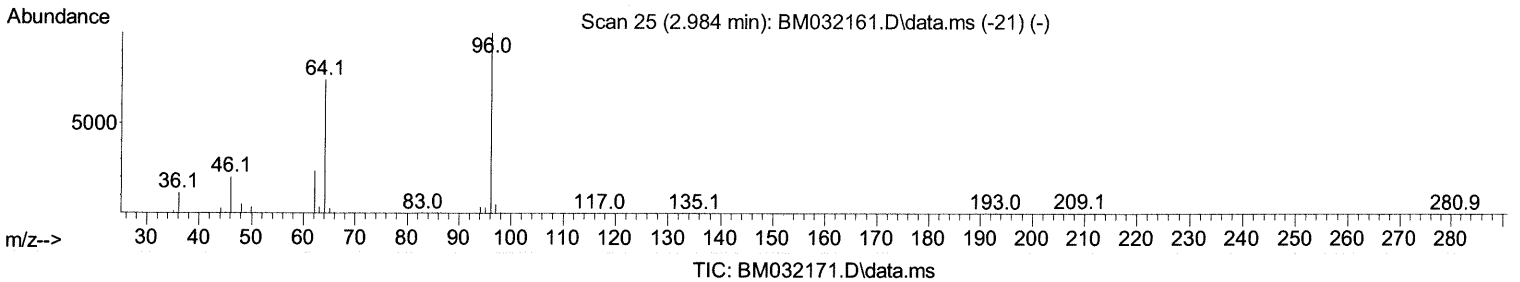
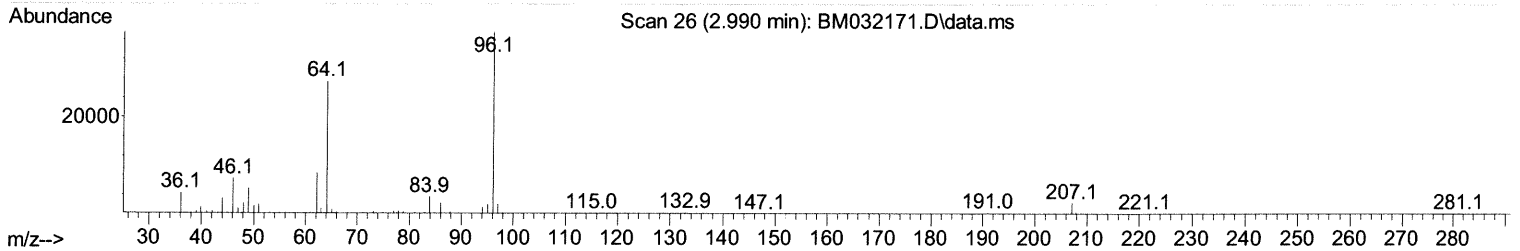
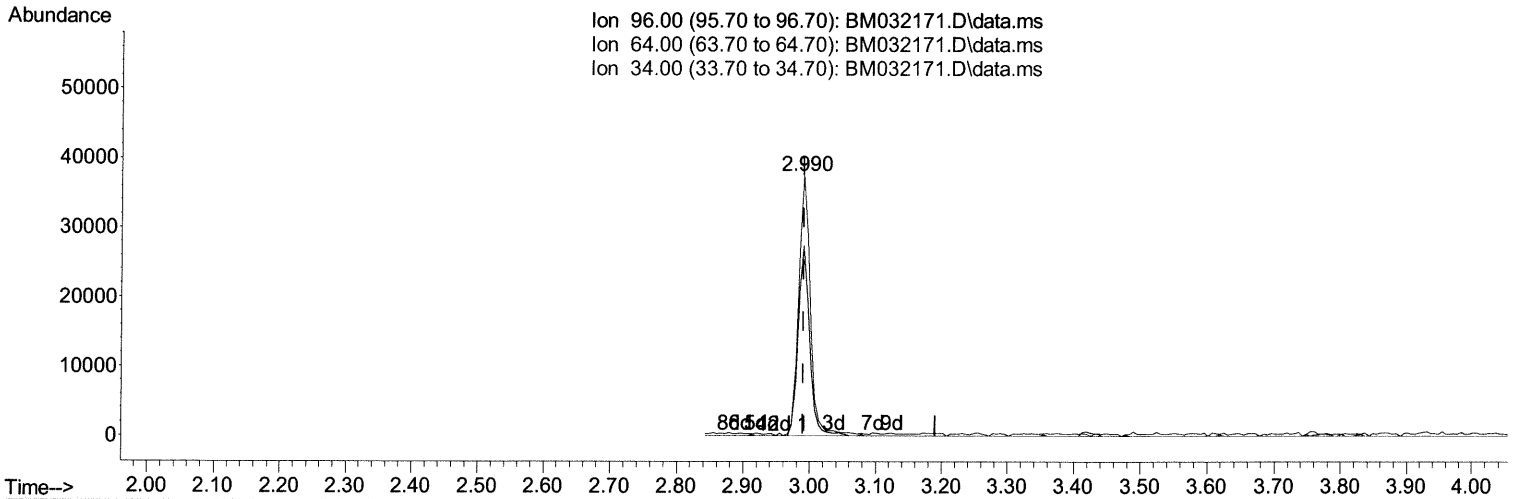
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM032171.D\data.ms

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

2.990min (-0.000) 4.93 ng/uL m 09/29/21JU

response 46467

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.40	73.02
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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Manual Integrations
 APPROVED

mohammad
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	398059	20.000	ng/ul	0.00
20) Naphthalene-d8	10.460	136	1573560	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.313	164	876383	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.036	188	1600688	20.000	ng/ul	0.00
79) Chrysene-d12	21.195	240	1339299	20.000	ng/ul	0.00
88) Perylene-d12	23.365	264	1330880	20.000	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.990	96	46467m	4.929	ng/ul	0.00	09/29/21 JU
4) Pyridine-d5	3.419	84	271887	10.159	ng/ul	0.00	
7) Phenol-d5	6.848	99	258937	8.195	ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth...	6.990	67	730285	38.438	ng/ul	0.00	
11) 2-Chlorophenol-d4	7.201	132	755205	30.786	ng/ul	0.00	
15) 4-Methylphenol-d8	8.389	113	479517	19.408	ng/ul	0.00	
21) Nitrobenzene-d5	8.819	128	446616	47.520	ng/ul	0.00	
24) 2-Nitrophenol-d4	9.542	143	450277	45.309	ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.083	165	812266	35.754	ng/ul	0.00	
31) 4-Chloroaniline-d4	10.589	131	1056434	34.406	ng/ul	0.00	
46) Dimethylphthalate-d6	13.713	166	2630406	46.196	ng/ul	0.00	
49) Acenaphthylene-d8	14.001	160	3262566	45.098	ng/ul	0.00	
54) 4-Nitrophenol-d4	14.507	143	63478	6.853	ng/ul	0.00	
60) Fluorene-d10	15.301	176	2292148	47.177	ng/ul	0.00	
65) 4,6-Dinitro-2-methylph...	15.413	200	299604	39.605	ng/ul	0.00	
73) Anthracene-d10	17.136	188	3547046	53.108	ng/ul	0.00	
81) Pyrene-d10	19.418	212	3656642	51.892	ng/ul	0.00	
92) Benzo(a)pyrene-d12	23.230	264	3276837	52.464	ng/ul	0.00	

Target Compounds

Qvalue

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