

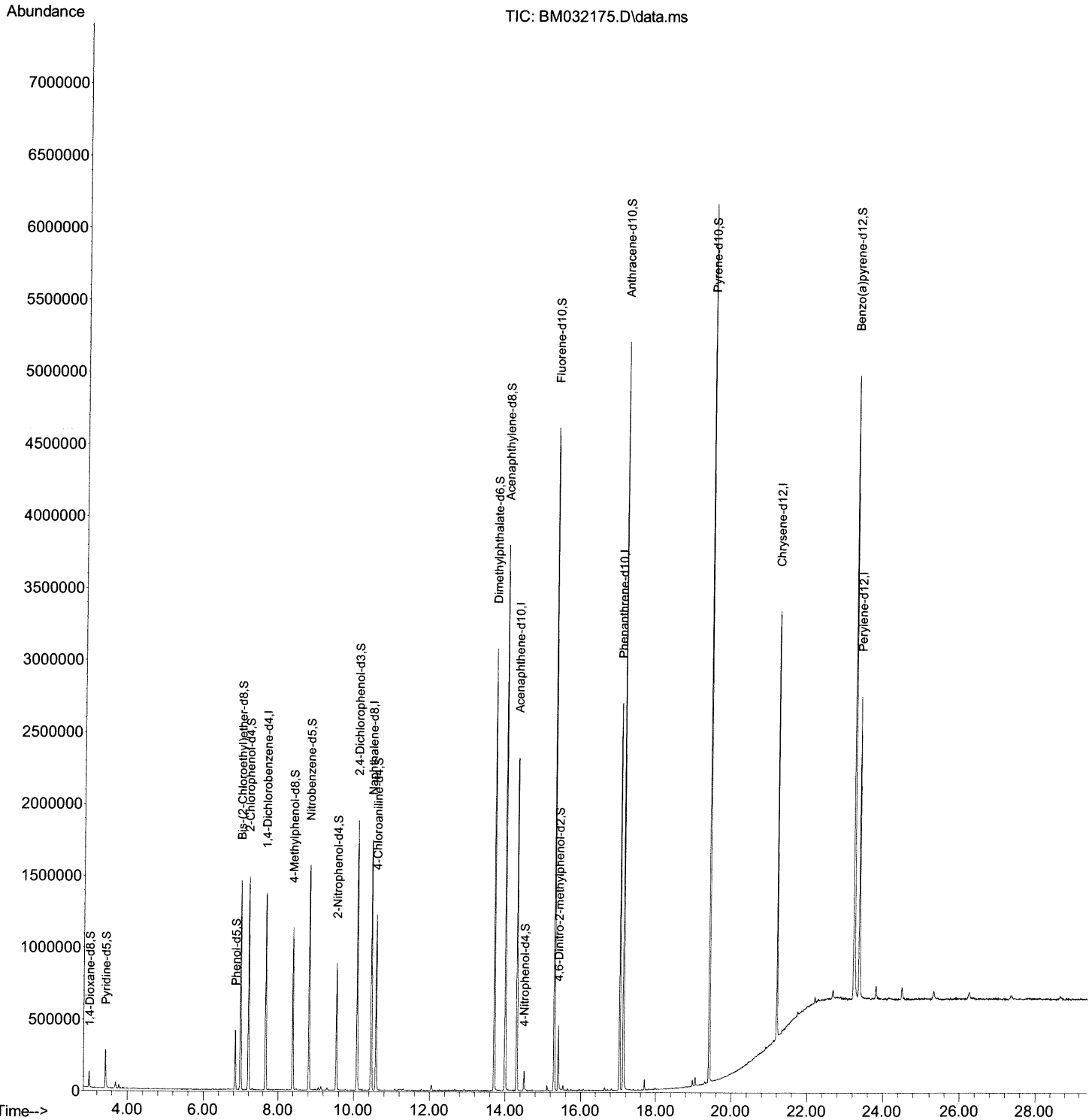
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
Data File : BM032175.D
Acq On : 26 Sep 2021 04:29
Operator : CG/JU
Sample : M3919-17
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GB6Z6

Manual Integrations
APPROVED

mohammad
9/27/2021 3:06:42 PM

Quant Time: Sep 26 05:07:30 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)

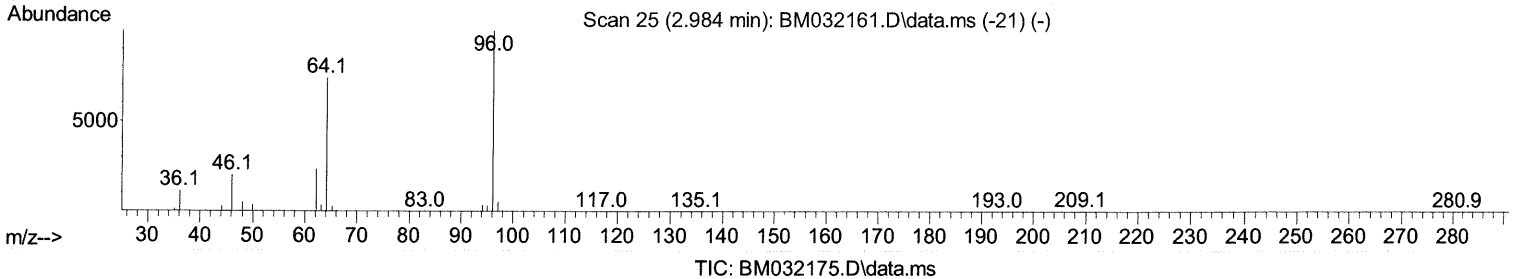
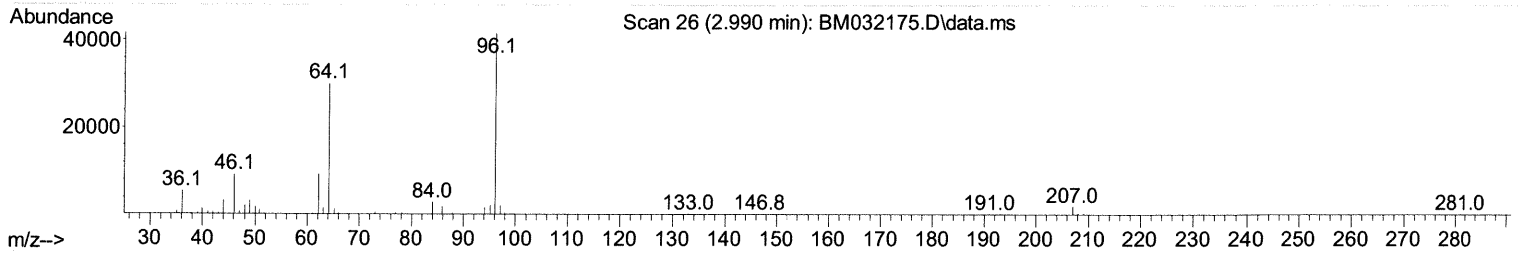
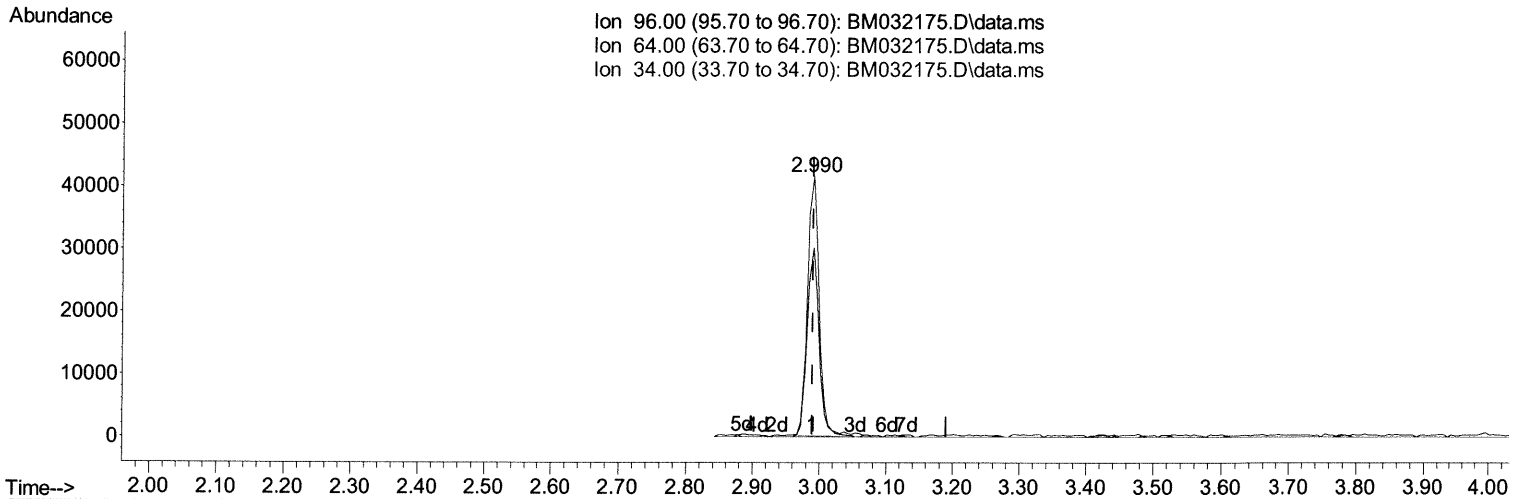
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TIC: BM032175.D\data.ms

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

2.990min (-0.000) 5.95 ng/uL

response 50987

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.40	72.43
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

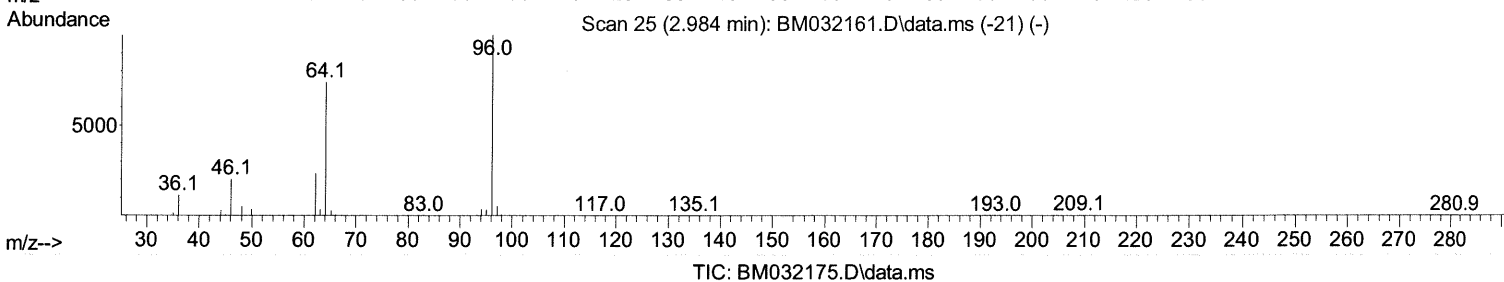
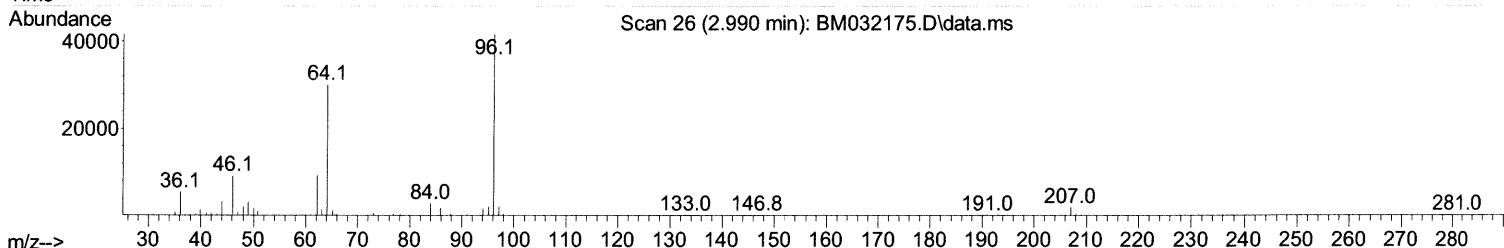
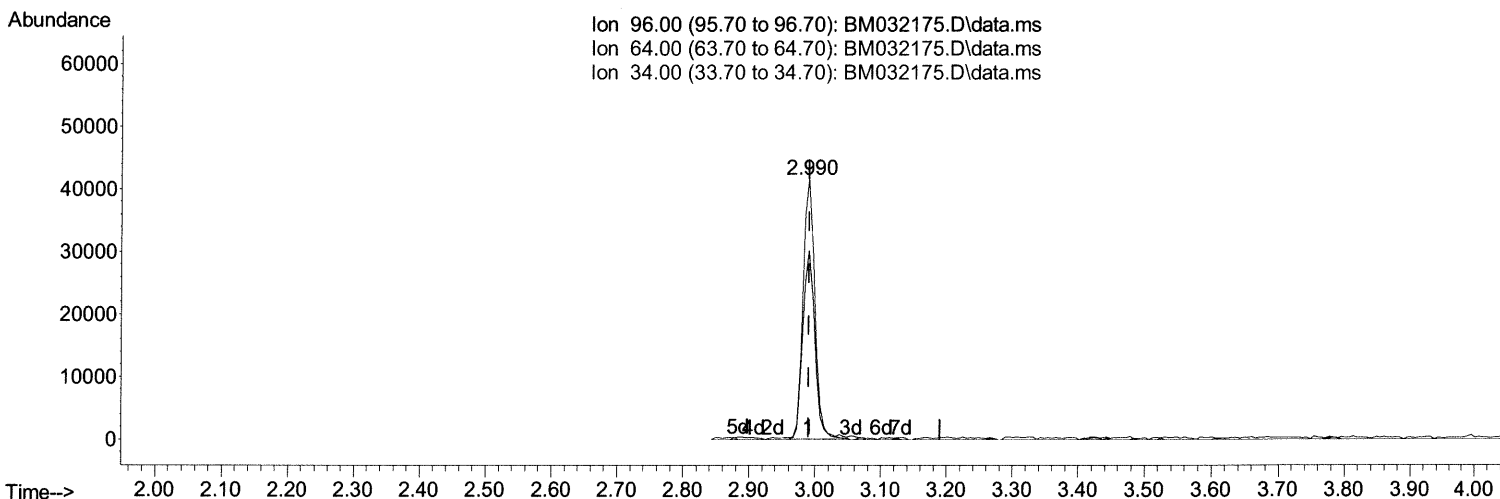
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TIC: BM032175.D\data.ms

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

2.990min (-0.000) 6.00 ng/uL m 09/29/21 JU

response 51404

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.40	72.43
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.666	152	362022	20.000 ng/ul	0.00
20) Naphthalene-d8	10.460	136	1389164	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.312	164	784604	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.036	188	1480406	20.000 ng/ul	0.00
79) Chrysene-d12	21.194	240	1321714	20.000 ng/ul	0.00
88) Perylene-d12	23.359	264	1321572	20.000 ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.990	96	51404m	> 5.995 ng/ul	>0.00	09/29/21JU
4) Pyridine-d5	3.419	84	138802	5.703 ng/ul	0.00	
7) Phenol-d5	6.842	99	221132	7.695 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth...	6.984	67	607168	35.139 ng/ul	0.00	
11) 2-Chlorophenol-d4	7.201	132	640349	28.703 ng/ul	0.00	
15) 4-Methylphenol-d8	8.383	113	397203	17.677 ng/ul	-0.01	
21) Nitrobenzene-d5	8.819	128	361196	43.532 ng/ul	0.00	
24) 2-Nitrophenol-d4	9.542	143	257702	29.373 ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.083	165	644448	32.132 ng/ul	0.00	
31) 4-Chloroaniline-d4	10.589	131	609115	22.471 ng/ul	0.00	
46) Dimethylphthalate-d6	13.713	166	1984907	38.937 ng/ul	0.00	
49) Acenaphthylene-d8	14.001	160	2480296	38.295 ng/ul	0.00	
54) 4-Nitrophenol-d4	14.501	143	26986	3.254 ng/ul	-0.01	
60) Fluorene-d10	15.295	176	1722929	39.609 ng/ul	0.00	
65) 4,6-Dinitro-2-methylph...	15.407	200	79915	11.422 ng/ul	-0.01	
73) Anthracene-d10	17.130	188	2788806	45.148 ng/ul	0.00	
81) Pyrene-d10	19.418	212	2982928	42.894 ng/ul	0.00	
92) Benzo(a)pyrene-d12	23.224	264	2735591	44.107 ng/ul	-0.01	

Target Compounds Qvalue

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