

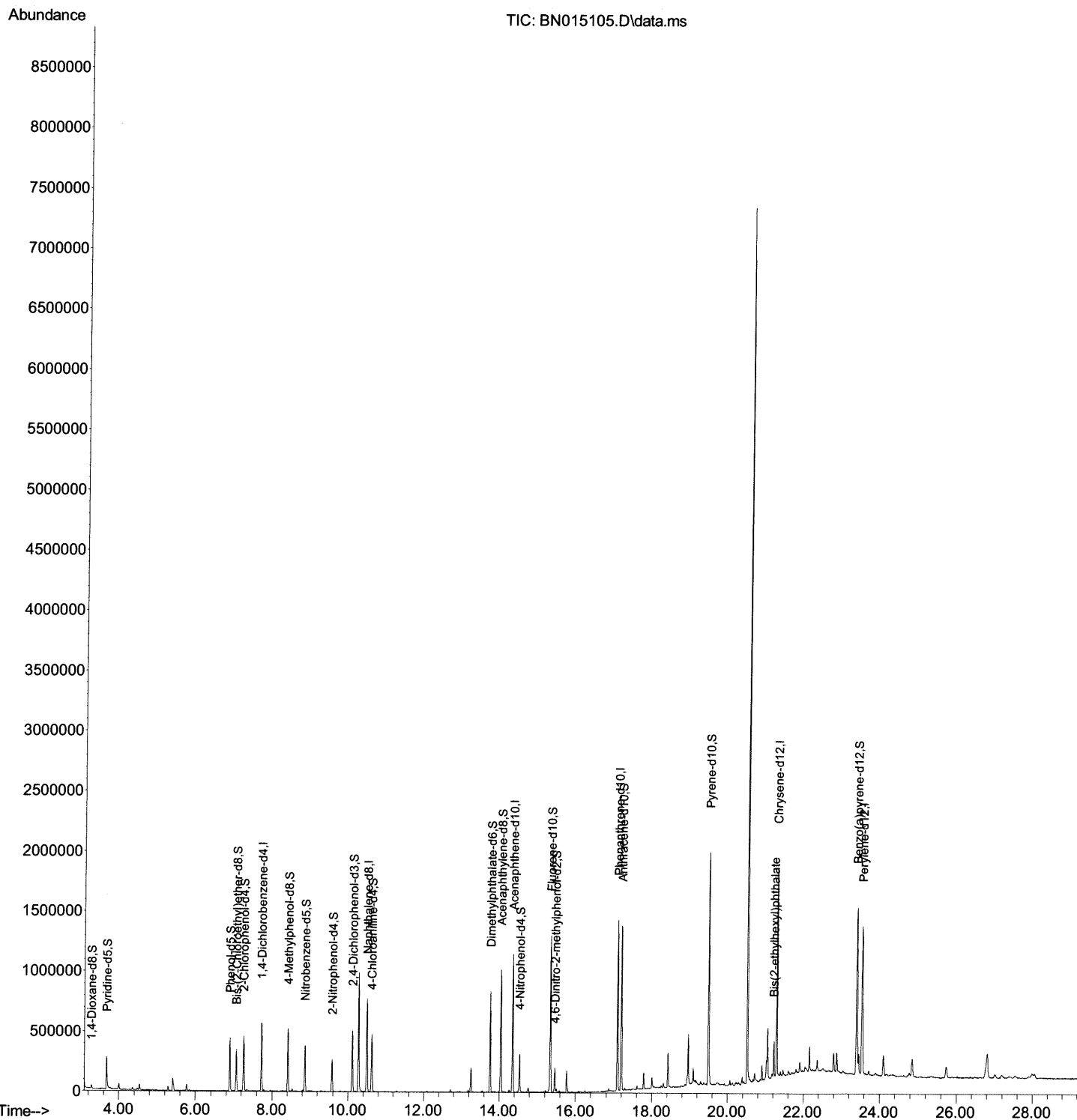
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB6

Manual Integrations
APPROVED

mohammad
6/28/2021 1:25:52 PM

Quant Time: Jun 26 23:58:33 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 24 14:43:13 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

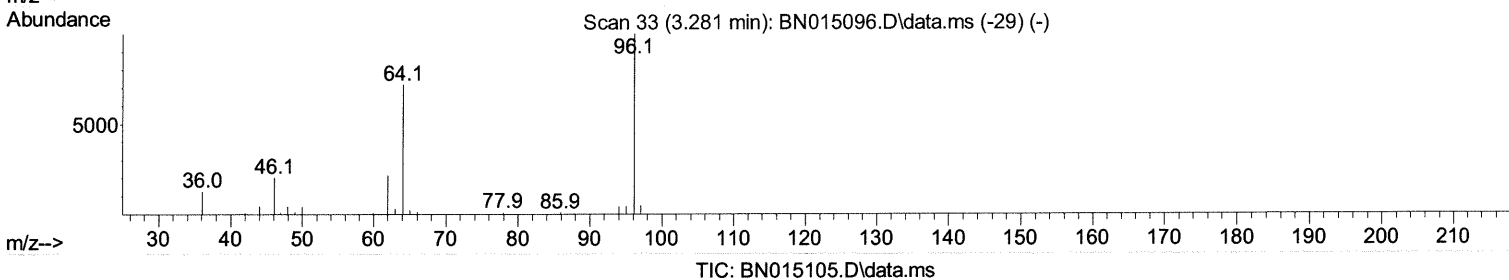
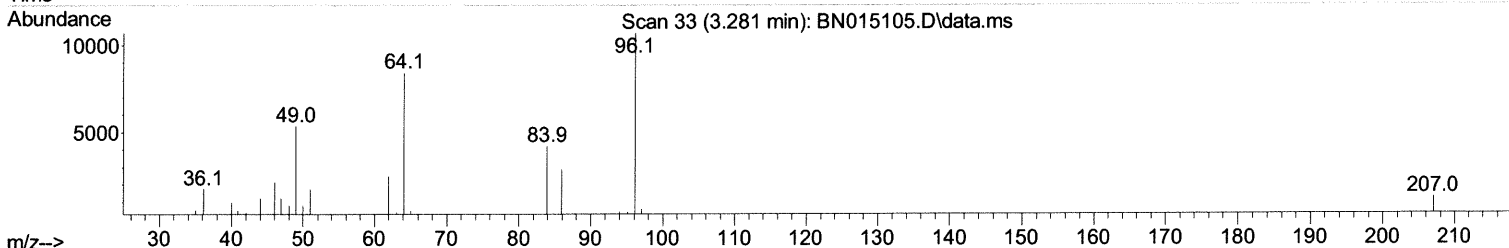
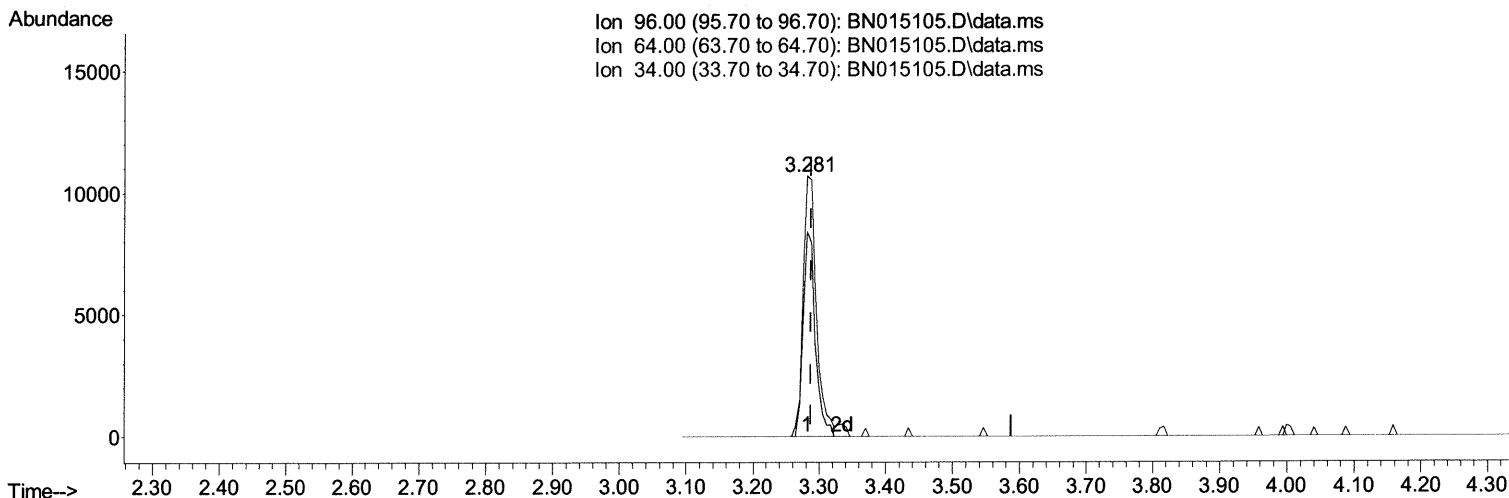
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(3) 1,4-Dioxane-d8 (S)

3.281min (-0.006) 3.78 ng/uL

response 15104

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.20	78.69
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

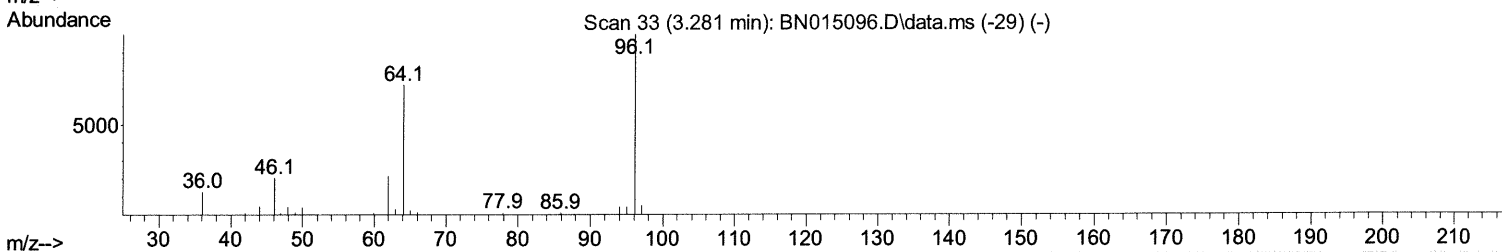
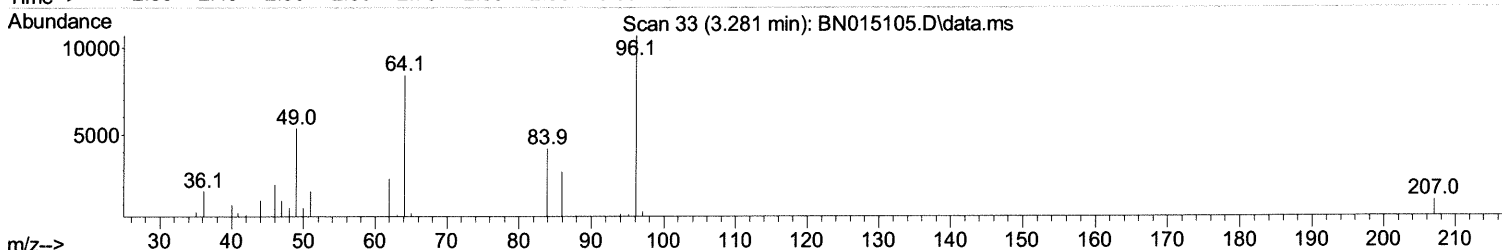
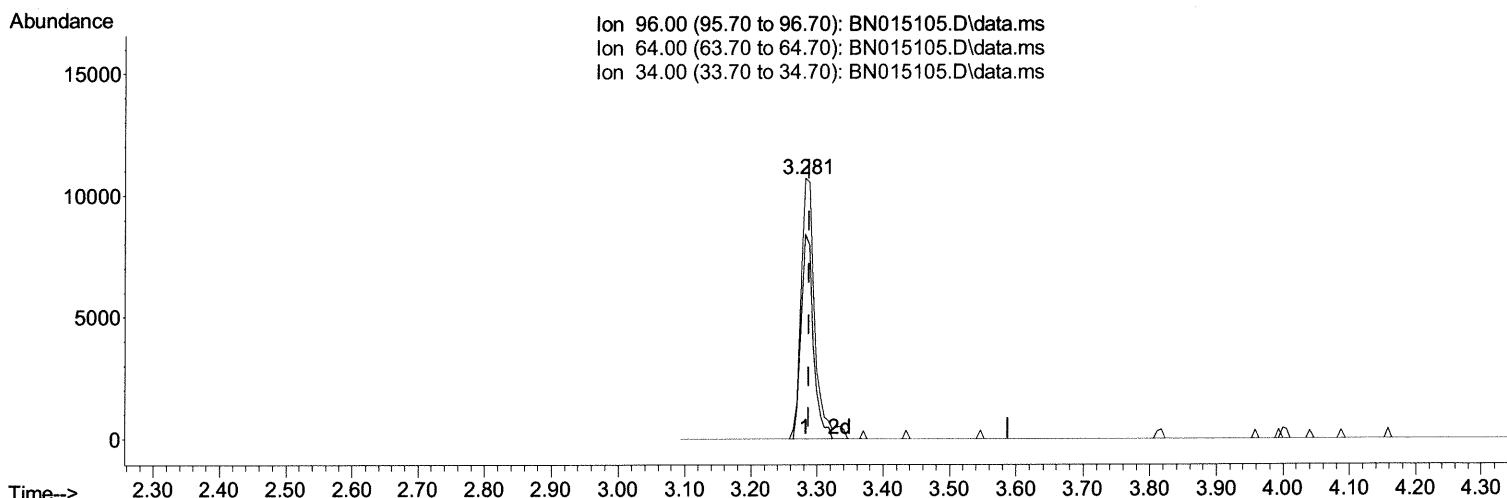
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Manual Integrations
APPROVED

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6/28/2021 1:25:52 PM

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BN015105.D\data.ms

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

3.281min (-0.006) 3.90 ng/uL m 06/29/2024

response 15564

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.20	78.69
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015105.D
 Acq On : 26 Jun 2021 08:02
 Operator : CG/JU
 Sample : M2704-07
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGDB6

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:25:52 PM

Quant Time: Jun 26 23:58:33 2021
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	155285	20.000	ng/ul	0.00
20) Naphthalene-d8	10.499	136	635606	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.351	164	386849	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	818489	20.000	ng/ul	0.00
79) Chrysene-d12	21.292	240	788805	20.000	ng/ul	0.00
88) Perylene-d12	23.539	264	818131	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	15564m	3.900	ng/ul	0.00
4) Pyridine-d5	3.676	84	150657	14.003	ng/ul	0.00
7) Phenol-d5	6.893	99	256607	19.039	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.064	67	150479	19.011	ng/ul	0.00
11) 2-Chlorophenol-d4	7.258	132	204477	20.008	ng/ul	0.00
15) 4-Methylphenol-d8	8.416	113	193697	18.621	ng/ul	0.00
21) Nitrobenzene-d5	8.869	128	93855	21.586	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	88019	22.727	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	181598	19.866	ng/ul	0.00
31) 4-Chloroaniline-d4	10.628	131	249383	17.603	ng/ul	0.00
46) Dimethylphthalate-d6	13.763	166	549361	20.681	ng/ul	0.00
49) Acenaphthylene-d8	14.040	160	705266	20.452	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	74422	16.062	ng/ul	0.00
60) Fluorene-d10	15.346	176	481960	20.855	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	37356	11.051	ng/ul	0.00
73) Anthracene-d10	17.192	188	773008	20.634	ng/ul	0.00
81) Pyrene-d10	19.498	212	938772	22.706	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.398	264	978620	24.230	ng/ul	0.00
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.222	149	91793	3.091	ng/ul	100

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