

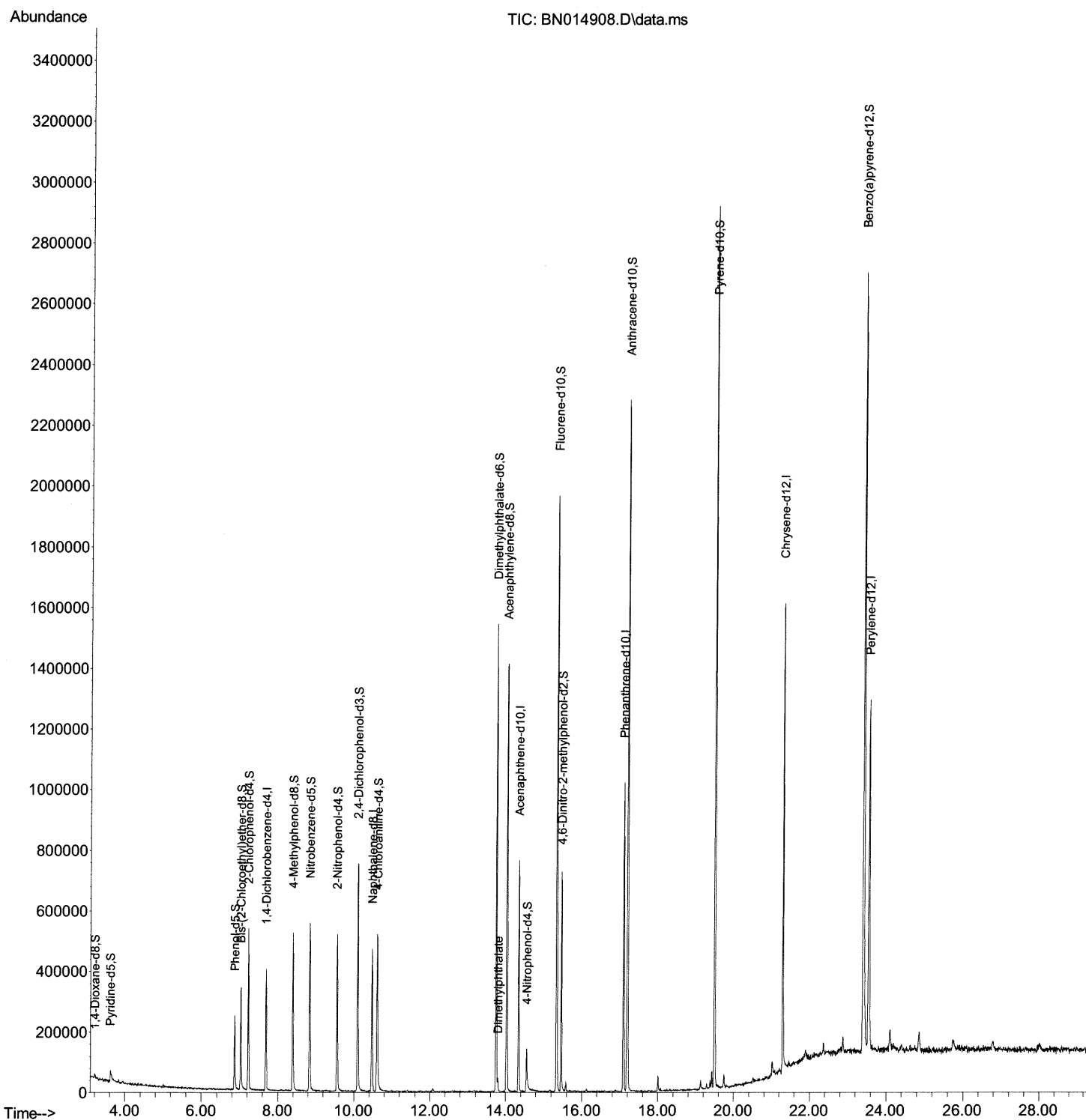
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061421\
Data File : BN014908.D
Acq On : 14 Jun 2021 21:03
Operator : CG/JU
Sample : M2649-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
B0AA7

Manual Integrations
APPROVED

mohammad
6/15/2021 2:42:13 PM

Quant Time: Jun 15 02:04:27 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061221MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 14 10:46:06 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

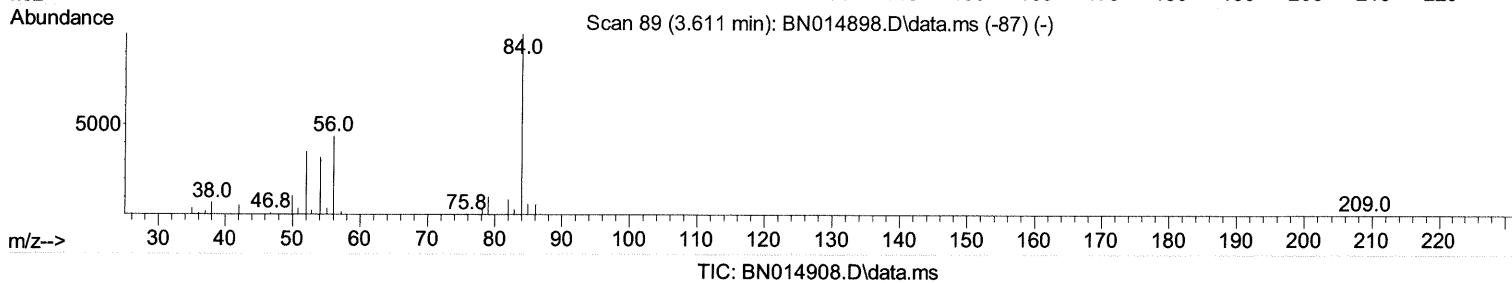
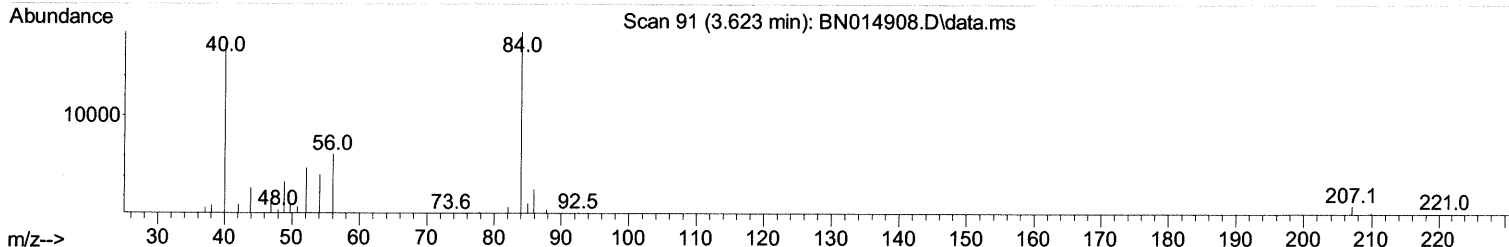
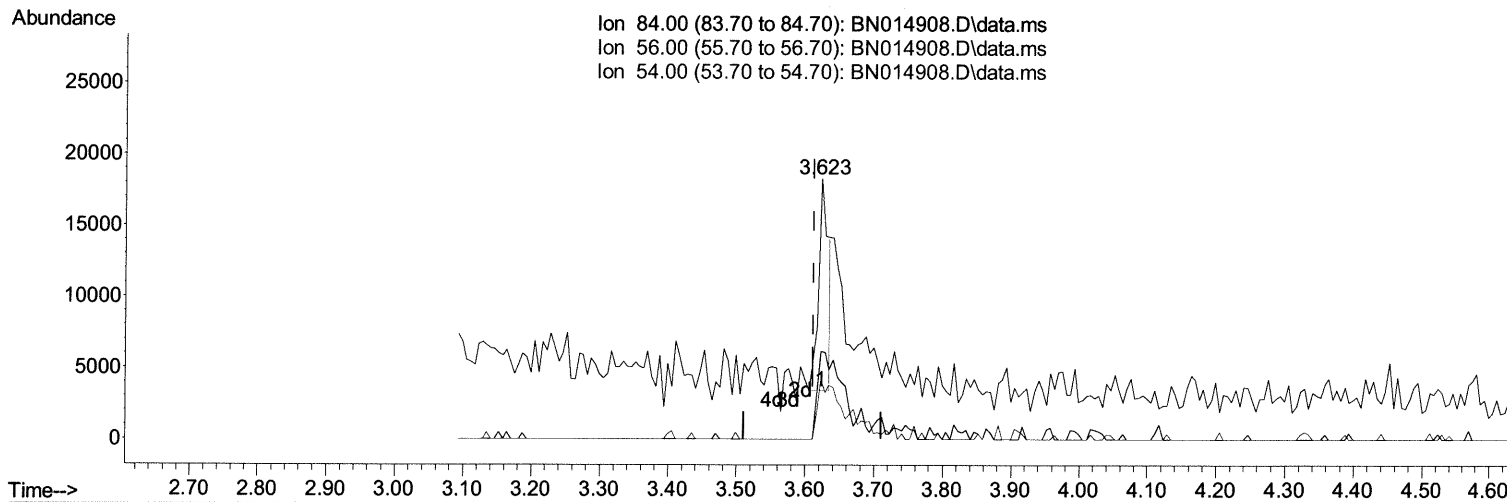
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(4) Pyridine-d5 (S)

3.623min (+ 0.012) 3.81 ng/ul

response 15190

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	44.70	33.71#
54.00	24.70	22.81
0.00	0.00	0.00

Quantitation Report (Qedit)

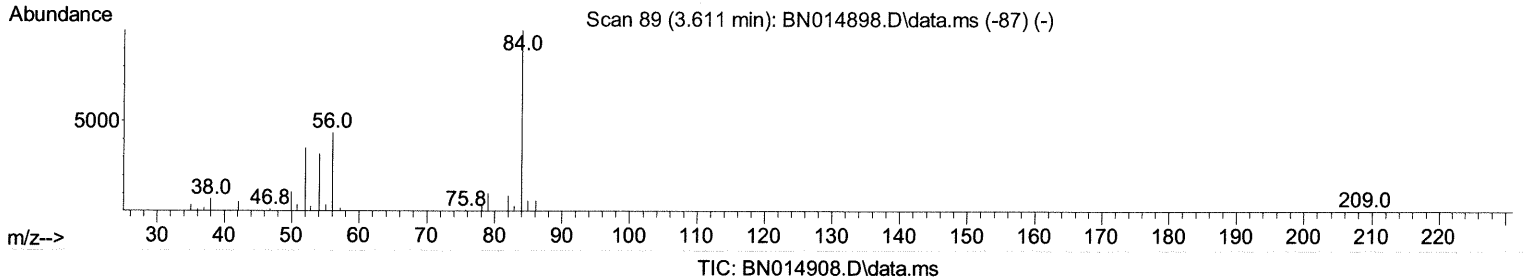
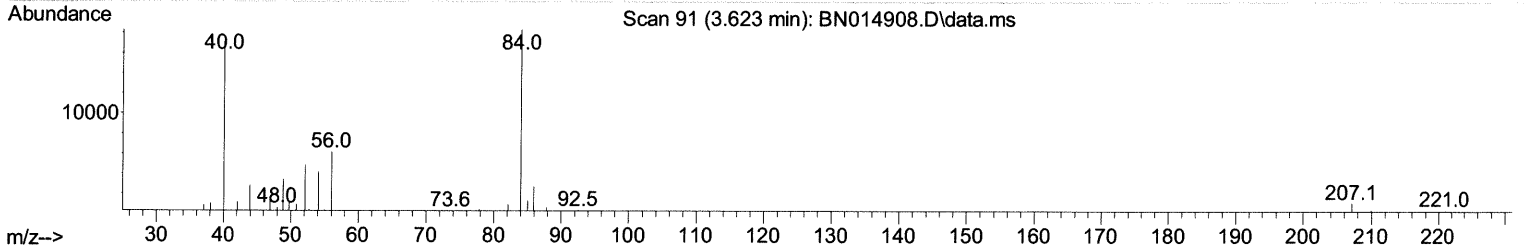
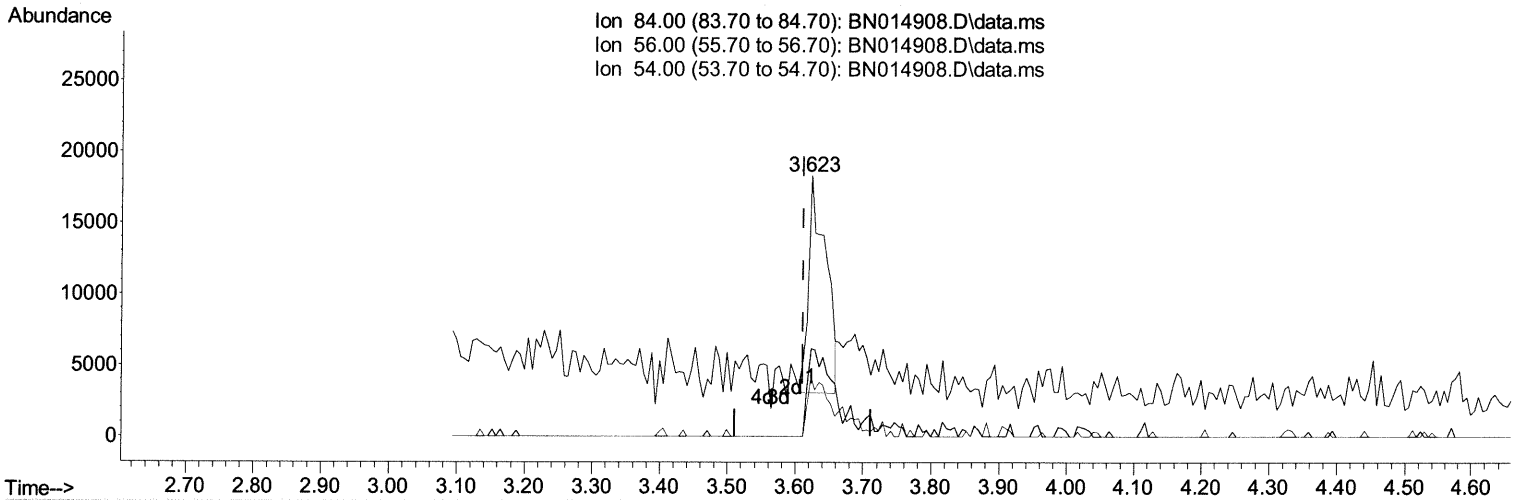
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Manual Integrations
 APPROVED

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 Response via : Initial Calibration



(4) Pyridine-d5 (S)

3.623min (+ 0.012) 6.74 ng/ul m 06/16/21 JU

response 26897

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	44.70	33.71#
54.00	24.70	22.81
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061421\
 Data File : BN014908.D
 Acq On : 14 Jun 2021 21:03
 Operator : CG/JU
 Sample : M2649-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 B0AA7

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:42:13 PM

Quant Time: Jun 15 02:04:27 2021
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 14 10:46:06 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	91218	20.000	ng/ul	0.00
20) Naphthalene-d8	10.475	136	334683	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.340	164	234114	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	575546	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.292	240	733270	20.000	ng/ul	0.00
88) Perylene-d12	23.545	264	811553	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	9018	5.257	ng/ul	0.00
4) Pyridine-d5	3.623	84	26897m	6.741	ng/ul	> 0.01 06/16/21JU
7) Phenol-d5	6.864	99	124751	18.258	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	137023	42.331	ng/ul	0.00
11) 2-Chlorophenol-d4	7.222	132	190684	36.398	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	162974	27.989	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	110077	43.831	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	118782	41.706	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	224688	39.182	ng/ul	0.00
31) 4-Chloroaniline-d4	10.610	131	222087	32.702	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	877946	47.172	ng/ul	0.00
49) Acenaphthylene-d8	14.034	160	945751	46.491	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	32247	12.334	ng/ul	0.00
60) Fluorene-d10	15.340	176	711095	46.837	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	134333	39.047	ng/ul	0.00
73) Anthracene-d10	17.192	188	1197914	45.991	ng/ul	0.00
81) Pyrene-d10	19.498	212	1514373	44.776	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	1958759	48.294	ng/ul	0.00
Target Compounds						
47) Dimethylphthalate	13.798	163	20436	1.165	ng/ul#	Qvalue 78

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