

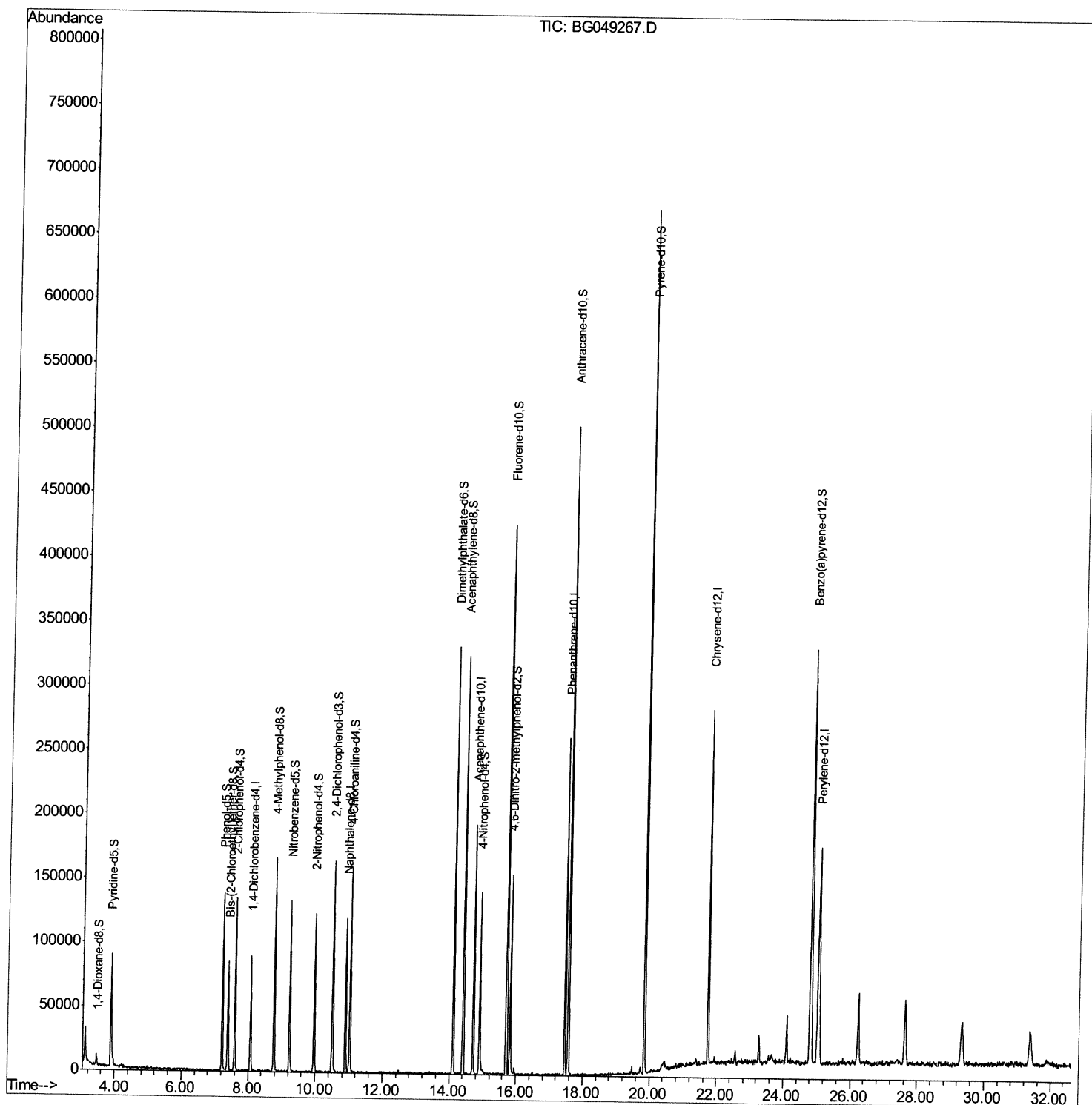
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
Data File : BG049267.D
Acq On : 10 Jul 2021 15:40
Operator : CG/JU
Sample : PB137573BL
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK573

Manual Integrations
APPROVED

mohammad
7/12/2021 12:34:47 PM

Quant Time: Jul 11 02:02:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 09 03:15:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

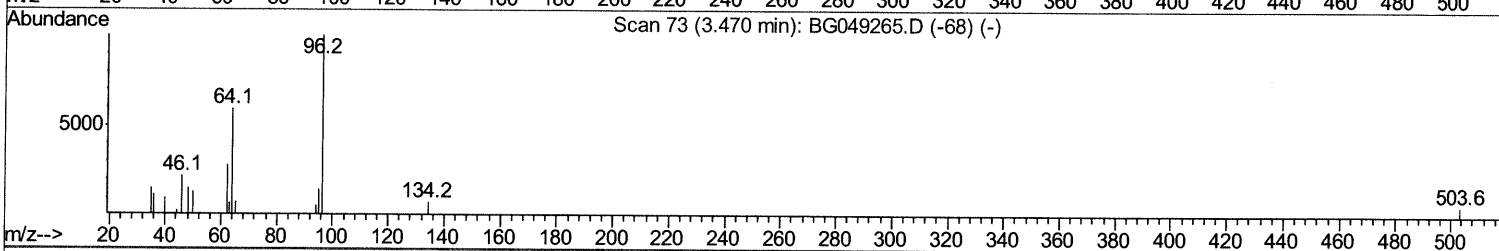
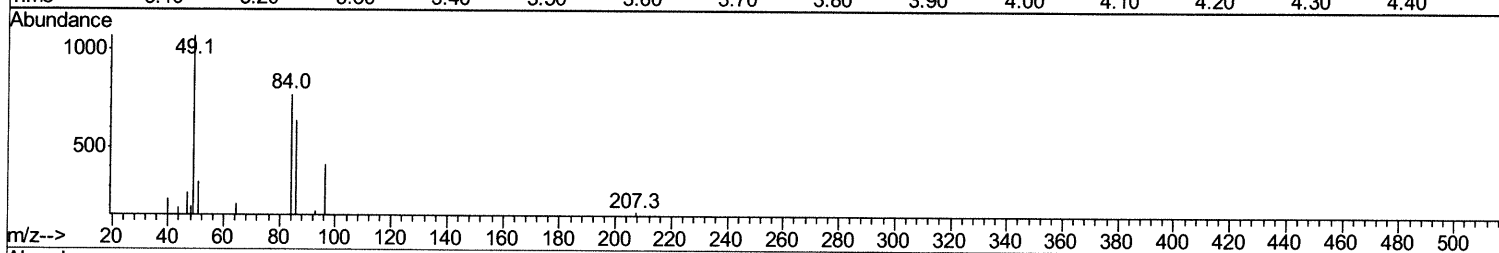
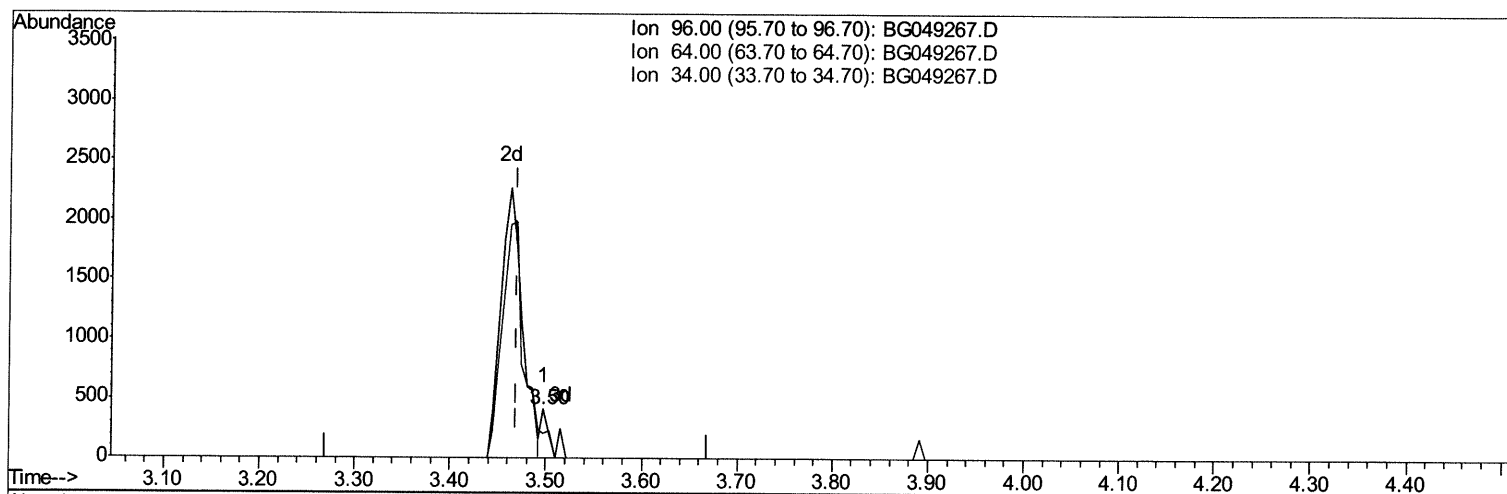
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TIC: BG049267.D

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

3.498min (+0.028) 0.46ng/uL

response 214

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	51.34
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

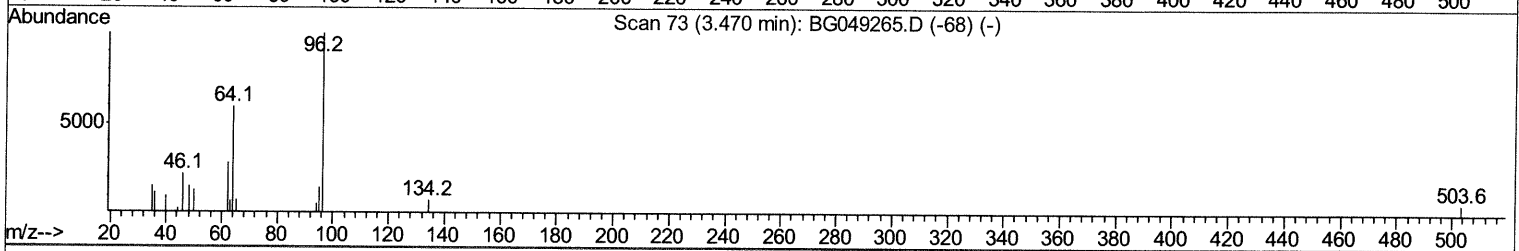
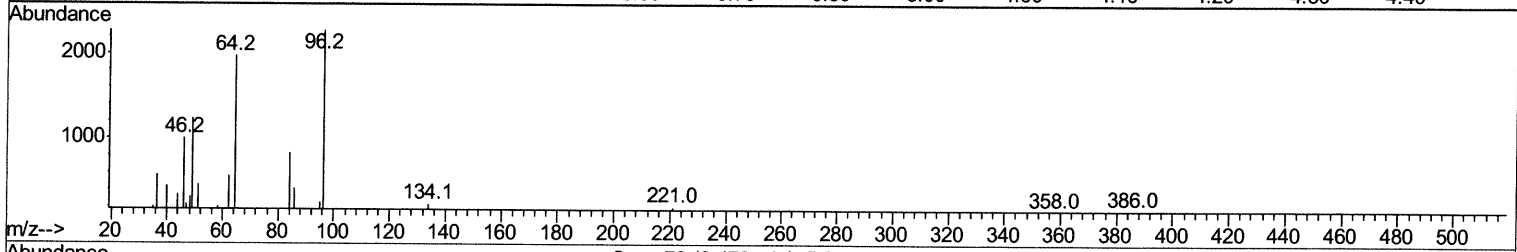
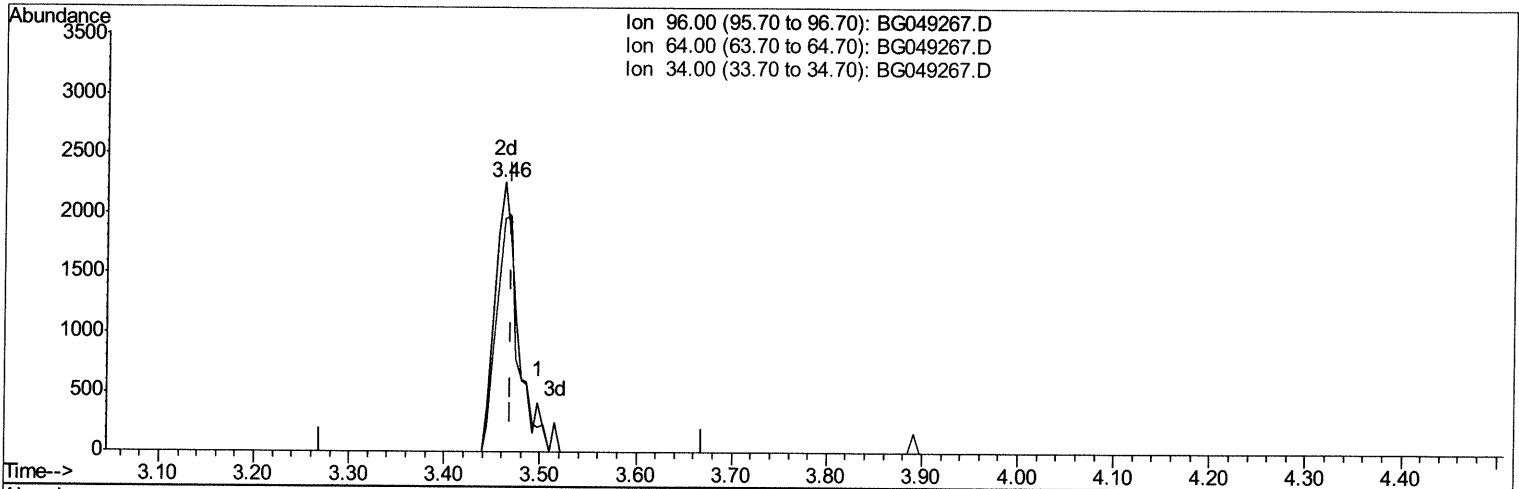
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Manual Integrations
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



TIC: BG049267.D

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

3.462min (-0.007) 7.99ng/uL m 07/13/21 JU

response 3705

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	86.59#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
 Data File : BG049267.D
 Acq On : 10 Jul 2021 15:40
 Operator : CG/JU
 Sample : PB137573BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK573

Manual Integrations
 APPROVED

mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	21817	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.89	136	96052	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	72386	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	170526	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	181516	20.00	ng/ul	0.00
88) Perylene-d12	25.04	264	181888	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	3705m>	7.99	ng/uL	> 0.00	07/13/21
4) Pyridine-d5	3.89	84	51290	37.61	ng/ul	0.00	
7) Phenol-d5	7.22	99	69789	36.77	ng/ul	0.00	
9) Bis-(2-Chloroethyl)ether-d	7.39	67	43520	39.54	ng/ul	0.00	
11) 2-Chlorophenol-d4	7.60	132	53880	38.41	ng/ul	0.00	
15) 4-Methylphenol-d8	8.77	113	57652	37.92	ng/ul	0.00	
21) Nitrobenzene-d5	9.24	128	27994	37.98	ng/ul	0.00	
24) 2-Nitrophenol-d4	9.97	143	31913	37.72	ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.51	165	58082	35.18	ng/ul	0.00	
31) 4-Chloroaniline-d4	11.02	131	77182	35.87	ng/ul	0.00	
46) Dimethylphthalate-d6	14.11	166	211126	37.93	ng/ul	0.00	
49) Acenaphthylene-d8	14.41	160	243550	37.29	ng/ul	0.00	
54) 4-Nitrophenol-d4	14.90	143	30916	33.55	ng/ul	-0.01	
60) Fluorene-d10	15.71	176	174185	36.96	ng/ul	0.00	
65) 4,6-Dinitro-2-methylphenol	15.82	200	33485	30.89	ng/ul	0.00	
73) Anthracene-d10	17.56	188	304988	39.28	ng/ul	0.00	
81) Pyrene-d10	19.85	212	378031	40.63	ng/ul	0.00	
92) Benzo(a)pyrene-d12	24.82	264	375019	39.66	ng/ul	0.00	

Target Compounds

Ovalue

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