

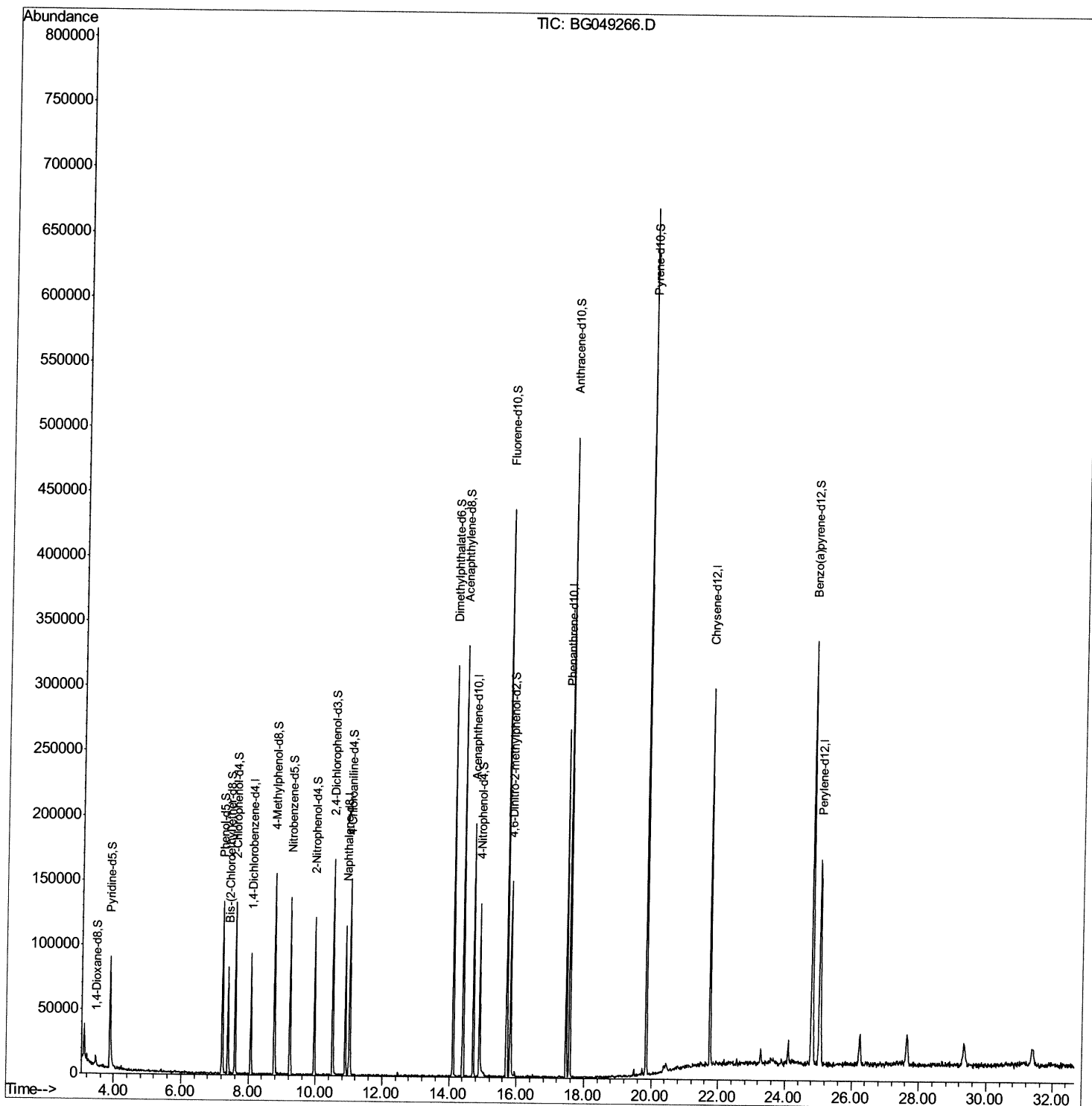
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
Data File : BG049266.D
Acq On : 10 Jul 2021 14:59
Operator : CG/JU
Sample : PB137517BL
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK517

Manual Integrations
APPROVED

mohammad
7/12/2021 12:35:16 PM

Quant Time: Jul 11 01:57:31 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)

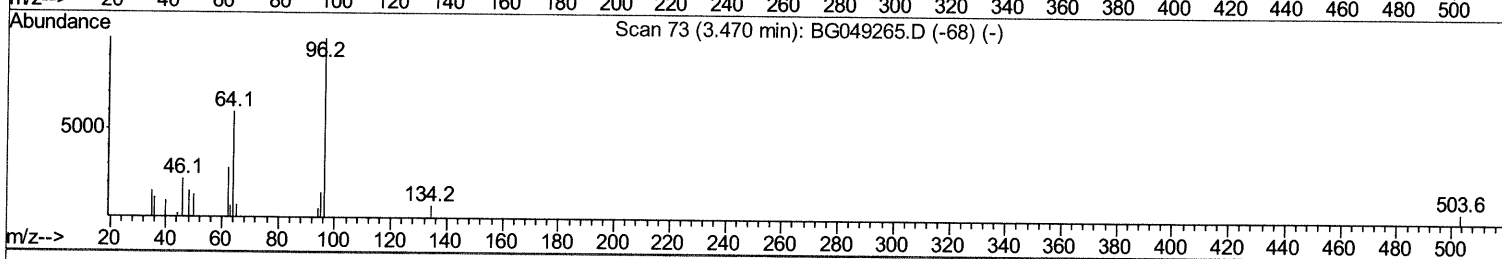
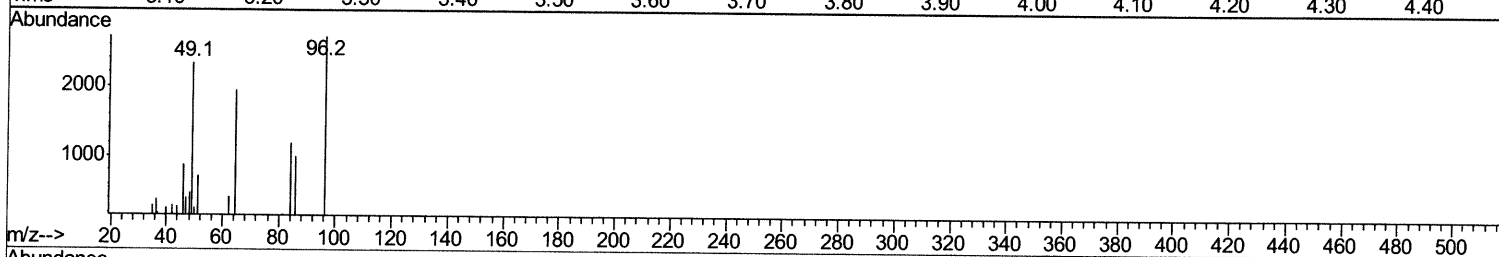
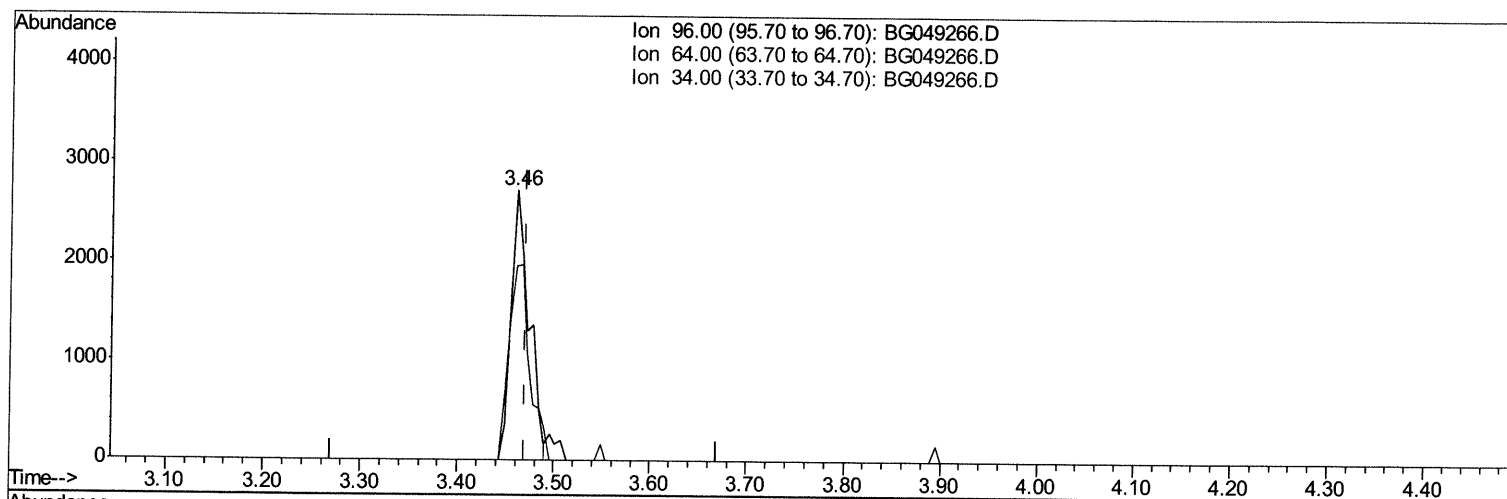
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TIC: BG049266.D

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

3.460min (-0.009) 7.13ng/uL

response 3511

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	71.67
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

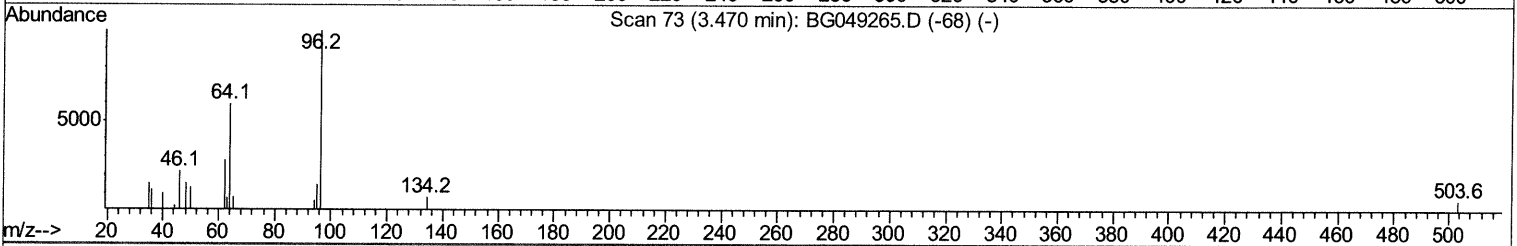
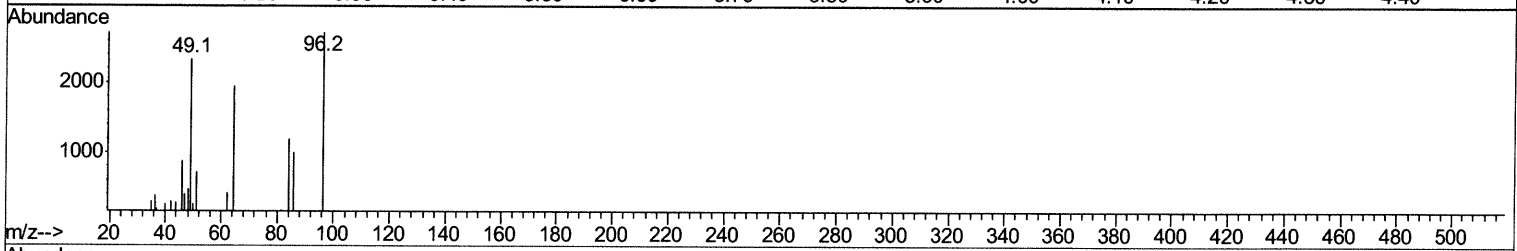
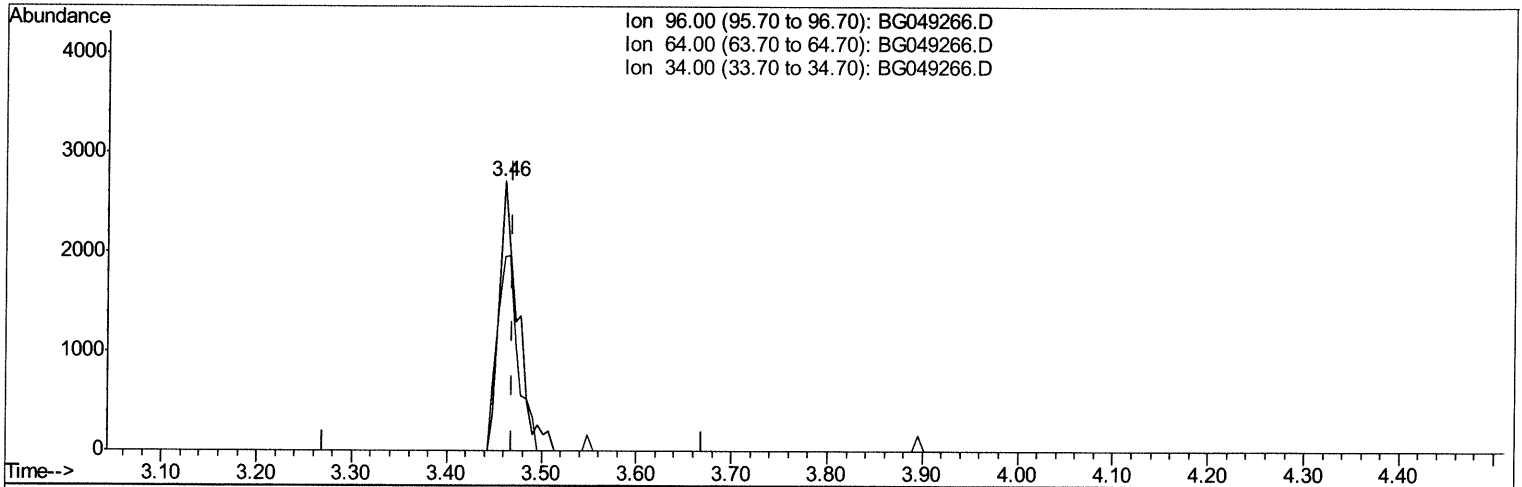
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Manual Integrations
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 Response via : Initial Calibration



TIC: BG049266.D

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

3.460min (-0.009) 7.58ng/uL m 07/13/21 JU

response 3731

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	71.67
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
 Data File : BG049266.D
 Acq On : 10 Jul 2021 14:59
 Operator : CG/JU
 Sample : PB137517BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
 APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	23148	20.00	ng/ul	0.00
20) Naphthalene-d8	10.89	136	97337	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	71180	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	169632	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	181414	20.00	ng/ul	0.00
88) Perylene-d12	25.04	264	183449	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	3731m >	7.58	ng/uL >	0.00	07/13/21 JU
4) Pyridine-d5	3.88	84	51111	35.32	ng/ul	0.00	
7) Phenol-d5	7.22	99	69980	34.75	ng/ul	0.00	
9) Bis-(2-Chloroethyl)ether-d	7.40	67	41355	35.41	ng/ul	0.00	
11) 2-Chlorophenol-d4	7.60	132	53192	35.74	ng/ul	0.00	
15) 4-Methylphenol-d8	8.77	113	56712	35.15	ng/ul	0.00	
21) Nitrobenzene-d5	9.24	128	28315	37.91	ng/ul	0.00	
24) 2-Nitrophenol-d4	9.96	143	32963	38.44	ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.51	165	59221	35.40	ng/ul	0.00	
31) 4-Chloroaniline-d4	11.03	131	76942	35.29	ng/ul	0.00	
46) Dimethylphthalate-d6	14.12	166	209079	38.19	ng/ul	0.00	
49) Acenaphthylene-d8	14.41	160	243124	37.85	ng/ul	0.00	
54) 4-Nitrophenol-d4	14.90	143	30696	33.88	ng/ul	-0.02	
60) Fluorene-d10	15.71	176	178620	38.54	ng/ul	0.00	
65) 4,6-Dinitro-2-methylphenol	15.82	200	33202	30.79	ng/ul	0.00	
73) Anthracene-d10	17.57	188	299337	38.76	ng/ul	0.00	
81) Pyrene-d10	19.85	212	371803	39.99	ng/ul	0.00	
92) Benzo(a)pyrene-d12	24.82	264	378294	39.67	ng/ul	0.00	

Target Compounds Ovalue

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