Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG113021\

Data File : BG051280.D

Acq On : 30 Nov 2021 14:58

Operator : CG/JU Sample : M4839-05

Misc

ALS Vial : 9 Sample Multiplier: 1

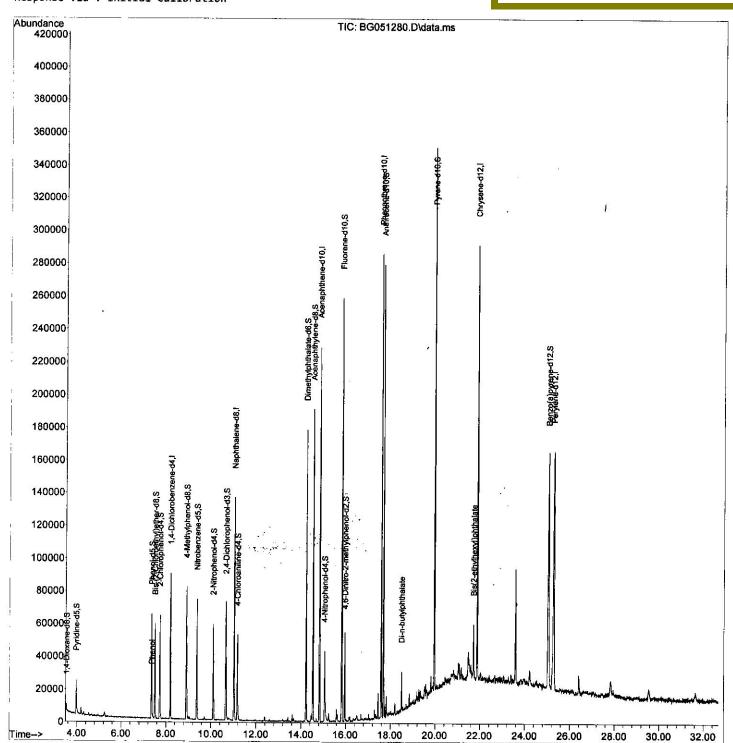
Quant Time: Nov 30 15:57:46 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : ESQP6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :mohammad ahmed 12/05/2021



SFAM-EPA-BG112321.M Tue Nov 30 16:00:58 2021

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG113021\

Data File : BG051280.D

Acq On : 30 Nov 2021 14:58

Operator : CG/JU Sample : M4839-05

Misc

ALS Vial : 9 Sample Multiplier: 1

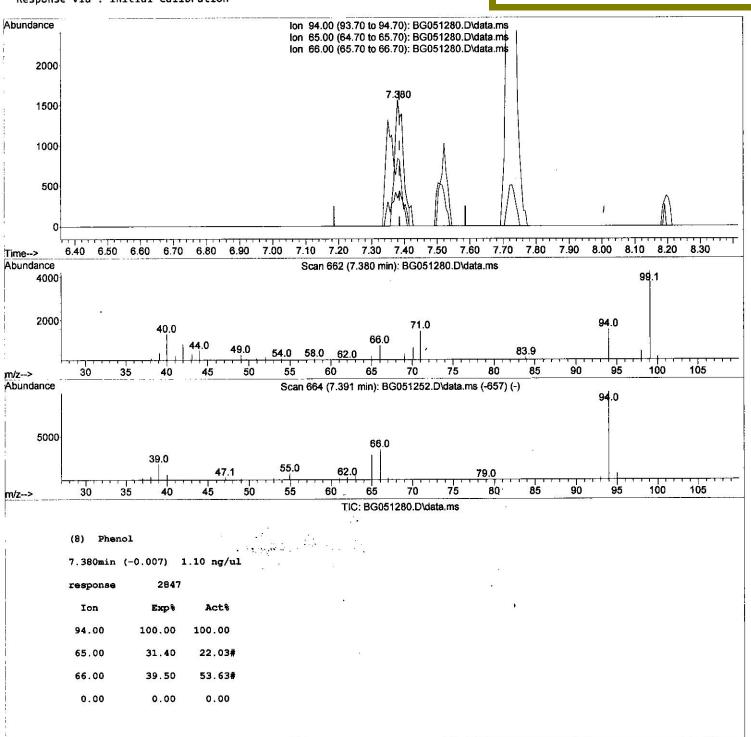
Quant Time: Nov 30 15:57:46 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument :
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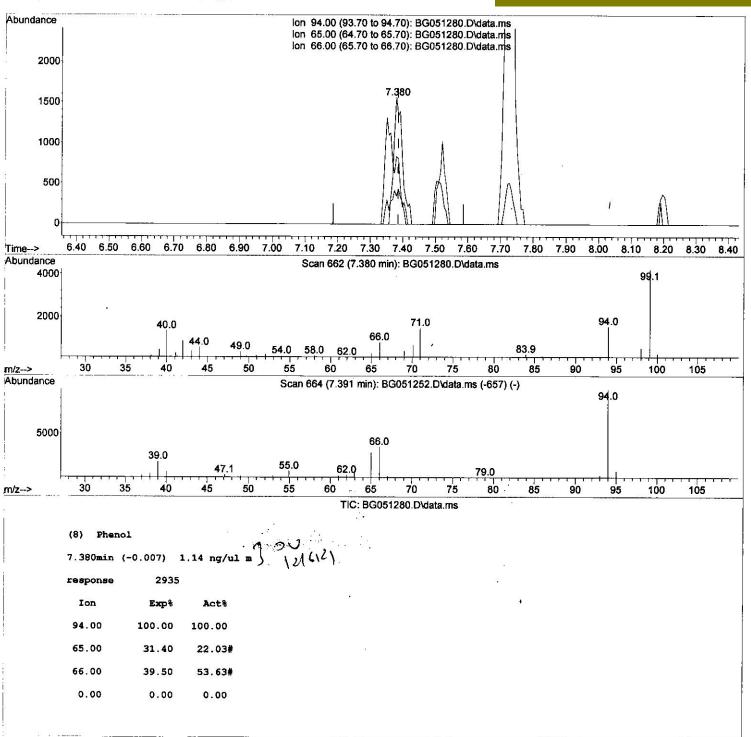
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Compound	R.T.	QIon	Response	Conc Units Dev(Min)	
Internal Standards					
 1,4-Dichlorobenzene-d4 	8.196	152	25220	20.000 ng/ul 0.0	0
20) Naphthalene-d8	11.023	136	116615	20.000 ng/ul 0.0	0
38) Acenaphthene-d10	14.824	164	81705	20.000 ng/ul 0.0	0
64) Phenanthrene-d10	17.574	188	176882	20.000 ng/ul 0.0	0
79) Chrysene-d12	21.875	240	158370	20.000 ng/ul 0.0	0
88) Perylene-d12	25.271	264	158153	20.000 ng/ul -0.0	1
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.531	96	2360	3.252 ng/uL -0.01	
4) Pyridine-d5	3.972	84	15153		200 20
7) Phenol-d5	7.356	99	41880	16.802 ng/ul 0.00	
9) Bis-(2-Chloroethyl)eth	7.509	67	29904	19.102 ng/ul 0.00	
11) 2-Chlorophenol-d4	7.726		31585	17.597 ng/ul 0.00	
15) 4-Methylphenol-d8	8.907	113	33798	16.803 ng/ul 0.00	
21) Nitrobenzene-d5	9.372	128	17546	17.824 ng/ul 0.00	
24) 2-Nitrophenol-d4	10.100	143	19518	17.577 ng/ul 0.00	
28) 2,4-Dichlorophenol-d3	10.652	165	31237	16.580 ng/ul 0.00	
31) 4-Chloroaniline-d4	11.164	131	32385	11.747 ng/ul 0.00	
46) Dimethylphthalate-d6	14.219	166	119252	18.969 ng/ul 0.00	
49) Acenaphthylene-d8	14.524	160	137419	17.335 ng/ul 0.00	
54) 4-Nitrophenol-d4	15.053		14148	13.903 ng/ul 0.00	
60) Fluorene-d10	15.817	176	102594	18.122 ng/ul 0.00	
65) 4,6-Dinitro-2-methylph	15.952	200	12397		
73) Anthracene-d10	17.674	188	163546	19.333 ng/ul 0.00	
81) Pyrene-d10	19.953	212	180009	18.785 ng/ul 0.00	
92) Benzo(a)pyrene-d12	25.035		144620	17.122 ng/ul 0.00	
Target Compounds			0	Qvalue	20
8) Phenol	7.380	94	2935m		121
78) Di-n-butylphthalate	18.508	149	17286	•	121
86) Bis(2-ethylhexyl)phtha			13069	1.897 ng/ul# 95	

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

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