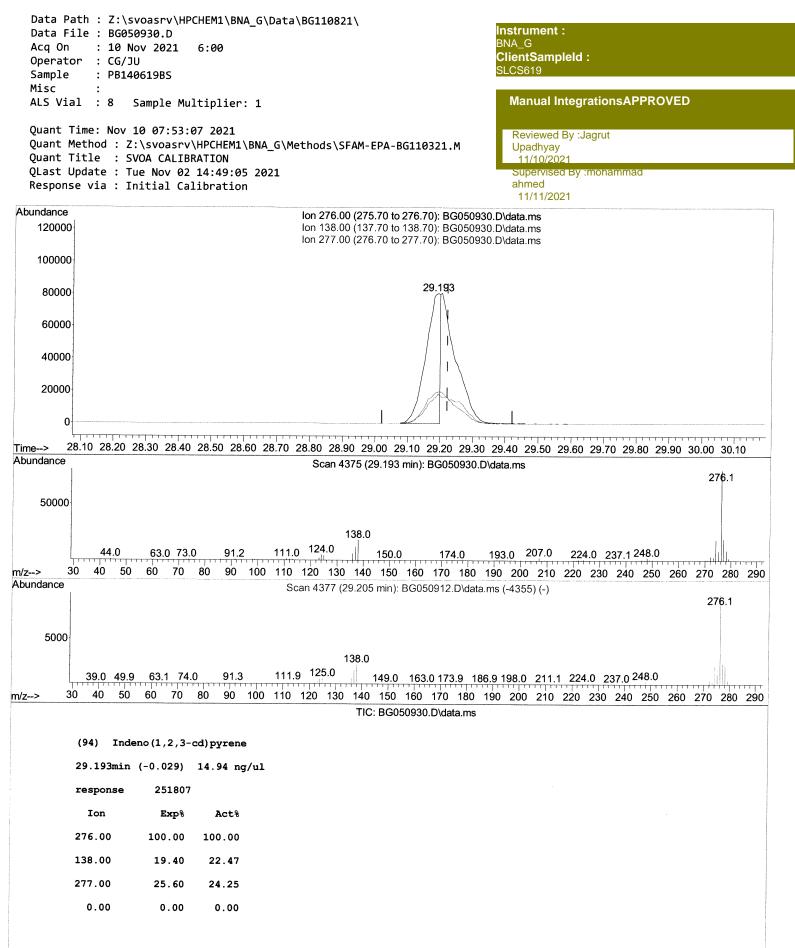
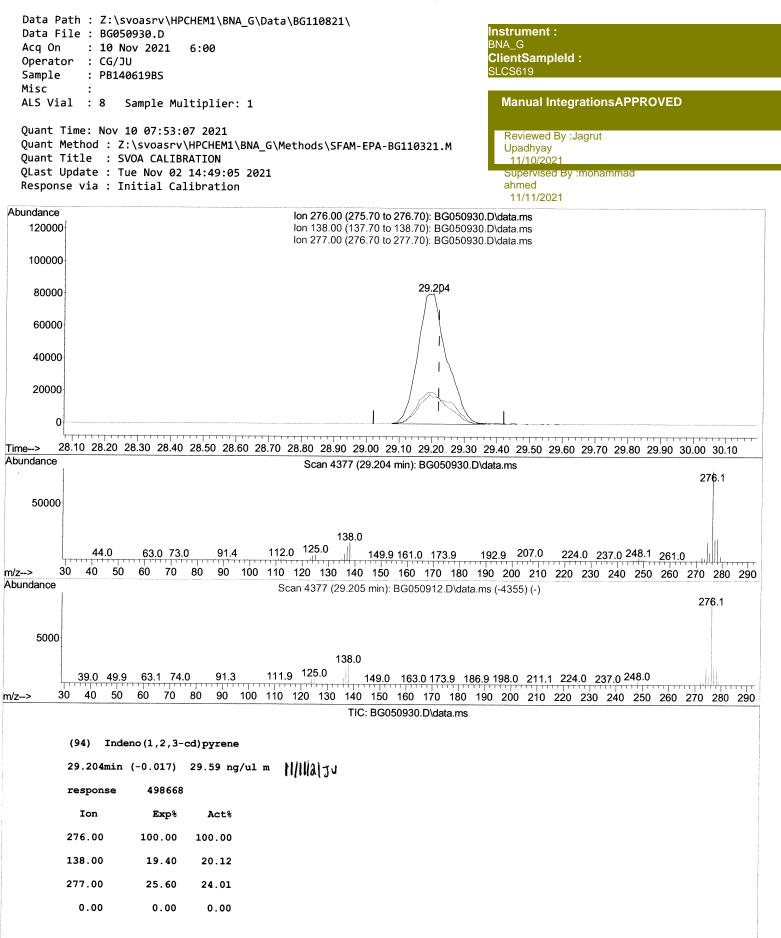


SFAM-EPA-BG110321.M Wed Nov 10 07:54:40 2021



Quantitation Report (Qedit)



Data Path : Z:\svoasrv\HPCHEM1	\BNA G\Da	nta\BG'	110821\				
Data File : BG050930.D	(500)_0(50		110021 (strument :	
Acq On : 10 Nov 2021 6:00						NA_G	
Operator : CG/JU						lientSampleId: _CS619	
Sample : PB140619BS Misc :					01		
ALS Vial : 8 Sample Multipl	ier· 1					Manual IntegrationsAPPROVED	
Quant Time: Nov 10 07:53:07 20					_	Reviewed By :Jagrut	
Quant Method : Z:\svoasrv\HPCH		\Metho	ods\SFAM-EF	PA-BG110321.M		Upadhyay	
Quant Title : SVOA CALIBRATIO QLast Update : Tue Nov 02 14:49						11/10/2021	
Response via : Initial Calibrat	105 2021					Supervised By :mohammad ahmed	20110712cm
						11/11/2021	100000
Compound	R.T.	QIon	Response	Conc Units De	ev(Min)		water and
							100000000000000000000000000000000000000
Internal Standards 1) 1,4-Dichlorobenzene-d4	0 220	450	44224				Contraction of Contra
20) Naphthalene-d8	8.229 11.055		44321 189584	20.00 ng/ul 20.00 ng/ul	0.00 0.00		Contractor (Contractor)
38) Acenaphthene-d10	14.851		118986	20.00 ng/ul	-0.01		inennen er
64) Phenanthrene-d10	17.595		258441	20.00 ng/ul	-0.01		COMPANY OF C
79) Chrysene-d12	21.895		222765	20.00 ng/ul	-0.01		unanesteres
88) Perylene-d12	25.291	264	223104	20.00 ng/ul	-0.02	2	Spannash
System Monitoring Compounds							SHOWER SHO
3) 1,4-Dioxane-d8	3.587	96	7571	5.51 ng/uL	0.00		SHIBUUU
4) Pyridine-d5	4.010	84	108742	26.47 ng/ul	0.00		oloninisco.
7) Phenol-d5	7.371		126034	26.66 ng/ul	0.00		1000000
9) Bis-(2-Chloroethyl)eth			82488	27.01 ng/ul	0.00		noninkis
11) 2-Chlorophenol-d4	7.753		87952	26.84 ng/ul	-0.01		SUSSIAN
15) 4-Methylphenol-d8 21) Nitrobenzene-d5	8.928 9.404		94505 45810	25.39 ng/ul	0.00		intrasicios
24) 2-Nitrophenol-d4	10.127		49802	28.43 ng/ul 27.80 ng/ul	0.00 -0.01		Streaming
28) 2,4-Dichlorophenol-d3	10.667		84872	28.13 ng/ul	0.00		Sanossia
31) 4-Chloroaniline-d4	11.184	131	107046	23.42 ng/ul	0.00		NAMERICAN
46) Dimethylphthalate-d6	14.246		260013	28.56 ng/ul	0.00		sittisenis
49) Acenaphthylene-d8 54) 4-Nitrophenol-d4	14.551		331327	29.21 ng/ul	0.00		vioriaiausia
60) Fluorene-d10	15.039 15.838		45666 229048	27.67 ng/ul 28.40 ng/ul	0.00 -0.01		and the day
65) 4,6-Dinitro-2-methylph			42779	27.30 ng/ul	0.00		(GRANNING)
73) Anthracene-d10	17.694		354159	28.98 ng/ul			and the second
81) Pyrene-d10	19.968		415339	28.87 ng/ul	-0.01		strassilles
92) Benzo(a)pyrene-d12	25.056	264	352459	28.58 ng/ul	-0.02		Subbisis
Target Compounds				0			unación de
2) 1,4-Dioxane	3.629	88	16584	ې 11.00 ng/uL	value 97		lainteente
5) Pyridine	4.034	79	113851	26.77 ng/ul	97		ilisii iniisiisii
6) Benzaldehyde	7.359	77	85102	28.54 ng/ul	94		STREET, STREET, ST
8) Phenol	7.401	94	130930	26.77 ng/ul	99		INTERACTOR INTERACTOR
 Bis(2-Chloroethyl)ether 2-Chlorophenol 	7.636	93 1 2 9	102323	27.95 ng/ul	98		annanna.
13) 2-Methylphenol	7.788 8.658	128 108	91728 95082	27.57 ng/ul 26.30 ng/ul	98 100		antoines
14) 2,2'-oxybis(1-Chloropr	8.746	45	153715	26.66 ng/ul	100 96		CERT/FILLOS
16) Acetophenone	9.057	105	149299	25.82 ng/ul	99		SAMORES
17) N-Nitroso-di-n-propyla	9.034	70	90783	26.02 ng/ul	96		anter anter a
18) 4-Methylphenol	8.993	108	101747	26.43 ng/ul	96		Constanting of the
19) Hexachloroethane 22) Nitrobenzene	9.316 9.445	117 77	39100 130850	28.11 ng/ul	96		antinging a
23) Isophorone	9.445	82	129850 244951	28.90 ng/ul 28.09 ng/ul	97 100		BUDIER
25) 2-Nitrophenol	10.156	139	53067	29.53 ng/ul	97		(1000 Land
26) 2,4-Dimethylphenol	10.203	107	115255	29.14 ng/ul	97		Billion.
27) Bis(2-Chloroethoxy)met	10.444	93	135685	28.88 ng/ul	98		OHGenter.
29) 2,4-Dichlorophenol	10.691	162	86223	29.29 ng/ul	97		-
30) Naphthalene 32) 4-Chloroaniline	11.102 11.208	128 127	298728 108767	28.82 ng/ul 23.97 ng/ul	97 100		
33) Hexachlorobutadiene	11.378	225	57119	29.57 ng/ul	100 98		Contraction of the local division of the loc
34) Caprolactam	11.966	113	32122	25.71 ng/ul	96		- contracting
35) 4-Chloro-3-methylphenol	12.312	107	104930	27.91 ng/ul	100		- Caraling
							È.

Data Path : Z:\svoasrv\HPCHEM1\	.BNA_G∖Da	ata\BG:	110821\		Instrument :
Data File : BG050930.D Acq On : 10 Nov 2021 6:00					BNA_G
Operator : CG/JU					ClientSampleId :
Sample : PB140619BS					SLCS619
Misc :					
ALS Vial : 8 Sample Multipli	er: 1				Manual IntegrationsAPPROVED
Quant Times New 10,07,52,07,200					
Quant Time: Nov 10 07:53:07 202 Quant Method : Z:\svoasrv\HPCHE		Mothe	de CEAM E	DA DC110331 M	Reviewed By :Jagrut
Quant Title : SVOA CALIBRATION		Metho	JUS \SFAM-E	PA-BG110321.M	Upadhyay 11/10/2021
QLast Update : Tue Nov 02 14:49					Supervised By :mohammad
Response via : Initial Calibrat	ion				ahmed
					11/11/2021
Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.694	1/12	196554	 27 92 ng/ul	06
37) 1-Methylnaphthalene	12.094		190334	27.83 ng/ul 27.87 ng/ul	96 97
39) 1,2,4,5-Tetrachloroben			105462	30.42 ng/ul	96
40) Hexachlorocyclopentadiene			47306	28.41 ng/ul#	97
<pre>41) 2,4,6-Trichlorophenol</pre>	13.288	196	69034	30.43 ng/ul	98
42) 2,4,5-Trichlorophenol	13.364		72689	29.84 ng/ul	100
43) 1,1'-Biphenyl	13.687		263689	30.32 ng/ul	100
44) 2-Chloronaphthalene 45) 2-Nitroaniline	13.734		205672	30.18 ng/ul	98
43) 2-Nicioaniine 47) Dimethylphthalate	13.934 14.293		80886 269113	29.87 ng/ul	98 99
48) 2,6-Dinitrotoluene	14.422		57902	29.57 ng/ul 30.39 ng/ul	100
50) Acenaphthylene	14.580		338987	29.82 ng/ul	99
51) 3-Nitroaniline	14.757		53909	27.36 ng/ul	91
52) Acenaphthene	14.915	153	223446	29.90 ng/ul	97
53) 2,4-Dinitrophenol	14.968		25634	24.39 ng/ul	87
55) 4-Nitrophenol	15.056		41803	27.62 ng/ul	95
56) Dibenzofuran 57) 2,4-Dinitrotoluene	15.250		316181	29.55 ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	15.209		80264 57542	29.52 ng/ul 30.06 ng/ul	100 97
59) Diethylphthalate	15.650		284195	29.17 ng/ul	99
61) Fluorene	15.897		248327	29.33 ng/ul	99
62) 4-Chlorophenyl-phenyle			129040	29.26 ng/ul	99
63) 4-Nitroaniline	15.914	138	57852	29.60 ng/ul	96
66) 4,6-Dinitro-2-methylph	15.973		43488	28.46 ng/ul#	97
67) N-Nitrosodiphenylamine	16.096		219636	30.40 ng/ul	97
 68) 4-Bromophenyl-phenylether 69) Hexachlorobenzene 	16.772 16.895		79020	30.74 ng/ul	96
70) Atrazine	17.030	284 200	81572 88202	30.86 ng/ul 28.79 ng/ul	95 98
71) Pentachlorophenol	17.242	266	33783	27.84 ng/ul	96
72) Phenanthrene	17.642	178	417171	30.24 ng/ul	99
74) Anthracene	17.730	178	412435	29.80 ng/ul	100
75) 1,2,3,4-Tetrachloroben	13.658	216	109784	31.20 ng/uL	95
76) Pentachlorobenzene	15.168	250	95116	29.17 ng/uL	97
77) Carbazole	18.000	167	388646	31.33 ng/ul	98
78) Di-n-butylphthalate 80) Fluoranthene	18.529 19.639	149 202	494819 519948	30.35 ng/ul 30.11 ng/ul	100 99
82) Pyrene	20.003	202	496949	29.45 ng/ul	100
83) Butylbenzylphthalate	20.867	149	219765	30.29 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.778	252	141418	26.07 ng/ul	98
85) Benzo(a)anthracene	21.872	228	461350	29.92 ng/ul	98
86) Bis(2-ethylhexyl)phtha	21.743	149	312283	29.99 ng/ul	100
87) Chrysene	21.942	228	440750	29.92 ng/ul	99
89) Di-n-octyl phthalate	23.012	149	532623	29.32 ng/ul	100
90) Benzo(b)fluoranthene 91) Benzo(k)fluoranthene	24.204 24.275	252 252	468364 434435	29.47 ng/ul	99
93) Benzo(a)pyrene	24.275	252	434435 441262	29.13 ng/ul 29.15 ng/ul	100 100
94) Indeno(1,2,3-cd)pyrene	29.204	276		> 29.59 ng/ul >	IVINALTU
95) Dibenzo(a,h)anthracene	29.263	278	419157	29.40 ng/ul	99
96) Benzo(g,h,i)perylene	30.427	276	414920	29.42 ng/ul	98

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

SFAM-EPA-BG110321.M Wed Nov 10 07:54:38 2021

2