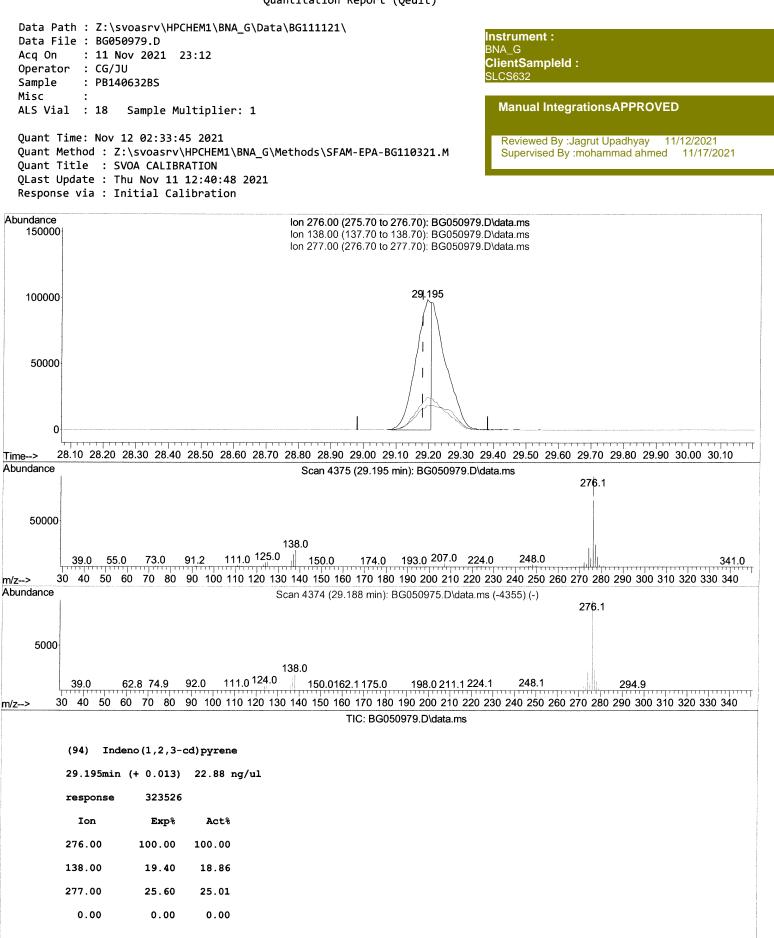
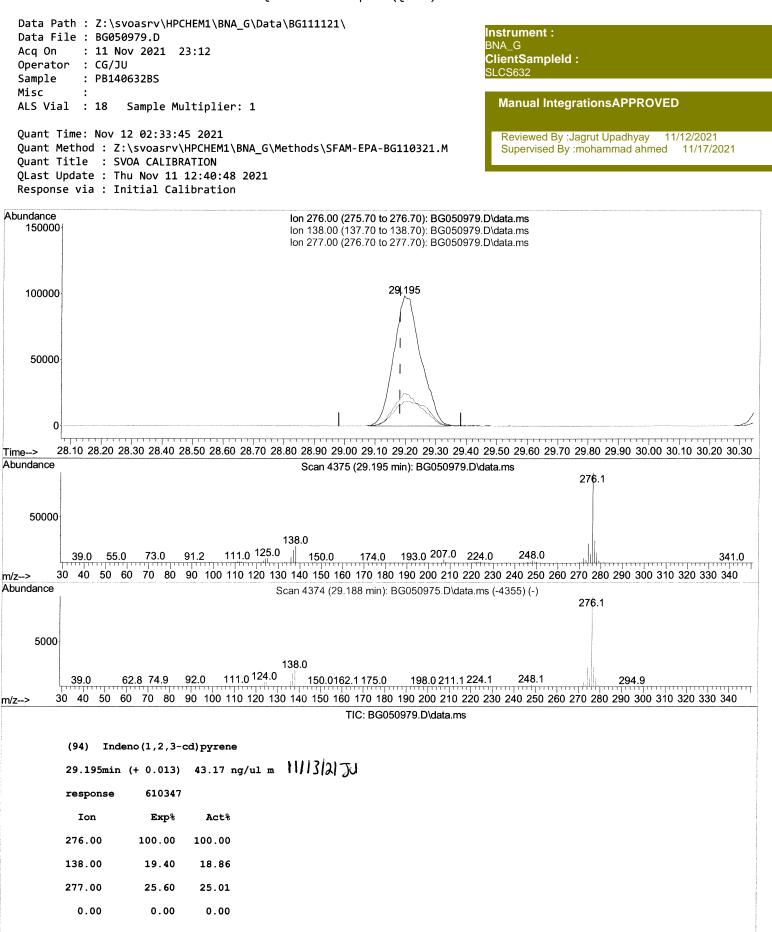


Quantitation Report (Qedit)



Quantitation Report (Qedit)



Data Path : Z:\svoasrv\HPCHEM1 Data File : BG050979.D Acq On : 11 Nov 2021 23:12 Operator : CG/JU Sample : PB140632BS Misc : ALS Vial : 18 Sample MultipJ Quant Time: Nov 12 02:33:45 202 Quant Method : Z:\svoasrv\HPCHE Quant Title : SVOA CALIBRATION QLast Update : Thu Nov 11 12:40 Response via : Initial Calibrat	_ 21 M1\BNA_G V 2:48 2021	Instrument : BNA_G ClientSampleId : SLCS632 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 11/12/2021 Supervised By :mohammad ahmed 11/17/2021				
Compound				Conc Units Dev		
Internal Standards						
1) 1,4-Dichlorobenzene-d4	0 771	150	24607	20.00 == (11	0.00	
, -	8.231		34607	20.00 ng/ul	0.00	
20) Naphthalene-d8	11.051		157964	20.00 ng/ul	0.00	
38) Acenaphthene-d10	14.853		104226	20.00 ng/ul	0.00	
64) Phenanthrene-d10	17.596		227085	20.00 ng/ul	0.00	
79) Chrysene-d12	21.891		190620	20.00 ng/ul	0.00	
88) Perylene-d12	25.293	264	187200	20.00 ng/ul	0.01	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	6174	5.76 ng/uL	0.00	
4) Pyridine-d5	4.018	84	96345	30.04 ng/ul	0.00	
7) Phenol-d5	7.373	99	133838	36.25 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth	7.543	67	83209	34.89 ng/ul	0.00	
<pre>11) 2-Chlorophenol-d4</pre>	7.755		93192	36.42 ng/ul	0.00	
15) 4-Methylphenol-d8	8.930		103294	35.54 ng/ul	0.00	
21) Nitrobenzene-d5	9.400	128	49524	36.89 ng/ul	0.00	
24) 2-Nitrophenol-d4	10.129	143	57045	38.22 ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.669	165	96272	38.29 ng/ul	0.00	
31) 4-Chloroaniline-d4	11.186	131	127090	33.38 ng/ul	0.00	
46) Dimethylphthalate-d6	14.247	166	320943	40.25 ng/ul	0.00	
49) Acenaphthylene-d8	14.547	160	392187	39.48 ng/ul	0.00	
54) 4-Nitrophenol-d4	15.046	143	62224	43.04 ng/ul	0.01	
60) Fluorene-d10	15.840	176	280712	39.74 ng/ul	0.00	
65) 4,6-Dinitro-2-methylph	15.957	200	60981	44.29 ng/ul	0.00	
73) Anthracene-d10	17.696	188	446333	41.57 ng/ul	0.00	
81) Pyrene-d10	19.970	212	512517	41.63 ng/ul	0.00	
92) Benzo(a)pyrene-d12	25.064	264	430817	41.63 ng/ul	0.01	
Target Compounds				Qv	value	
2) 1,4-Dioxane	3.630	88	14150	12.01 ng/uL	92	
5) Pyridine	4.036	79	106596	32.10 ng/ul	98	
6) Benzaldehyde	7.361	77	88421	37.97 ng/ul	97	
8) Phenol	7.402	94	143060	37.46 ng/ul	99	
10) Bis(2-Chloroethyl)ether	7.637	93	104260	36.47 ng/ul	100	
<pre>12) 2-Chlorophenol</pre>	7.790	128	98871	38.06 ng/ul	97	
<pre>13) 2-Methylphenol</pre>	8.666	108	105364	37.32 ng/ul	94	
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.754	45	163009	36.20 ng/ul	100	
16) Acetophenone	9.059	105	167161	37.02 ng/ul	99	
17) N-Nitroso-di-n-propyla	9.036	70	101321	37.19 ng/ul	99	
18) 4-Methylphenol	8.995	108	112729	37.51 ng/ul	99	
19) Hexachloroethane	9.318	117	38908	35.82 ng/ul	99	
22) Nitrobenzene	9.447	77	140811	37.61 ng/ul	96	
23) Isophorone	9.964	82	281764	38.78 ng/ul	99	
25) 2-Nitrophenol	10.158	139	61123	40.82 ng/ul	95	
26) 2,4-Dimethylphenol	10.205	107	126435	38.36 ng/ul	99	
27) Bis(2-Chloroethoxy)met	10.440	93	148382	37.90 ng/ul	98	
29) 2,4-Dichlorophenol	10.693	162	99145	40.43 ng/ul	96	
30) Naphthalene	11.104	128	332657	38.51 ng/ul	98	
32) 4-Chloroaniline	11.210	127	132160	34.96 ng/ul	98	
33) Hexachlorobutadiene	11.374	225	63203	39.26 ng/ul	97	
34) Caprolactam	11.979	113		• 39.28 ng/ul ≻		
35) 4-Chloro-3-methylphenol	12.314	107	125443	40.04 ng/ul	100	

Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG050979.D Acq On : 11 Nov 2021 23:12 Operator : CG/JU Sample : PB140632BS Misc : ALS Vial : 18 Sample Multipl: Quant Time: Nov 12 02:33:45 202:	BI C	NA_G IlentSampleId : LCS632 Manual IntegrationsAPPROVED				
Quant Method : Z:\svoasrv\HPCHEM Quant Title : SVOA CALIBRATION QLast Update : Thu Nov 11 12:40 Response via : Initial Calibrat	Reviewed By :Jagrut Upadhyay 11/12/2021 Supervised By :mohammad ahmed 11/17/2021					
Compound	R.T.	QIon	Response	Conc Units Dev	(Min)	
26) 2 Mothylpaphthalana	12 606	147				
36) 2-Methylnaphthalene	12.696		225192	38.27 ng/ul	98	
37) 1-Methylnaphthalene	12.914	142	228098	38.25 ng/ul	99	
39) 1,2,4,5-Tetrachloroben40) Hexachlorocyclopentadiene	13.055	216	127630	42.03 ng/ul	97	
40) Hexacillorocyclopentadiene 41) 2,4,6-Trichlorophenol	13.025	237	52580	36.05 ng/ul	98 100	
42) 2,4,5-Trichlorophenol	13.290 13.366	196 196	86993 94418	43.78 ng/ul	100 99	
43) 1,1'-Biphenyl	13.689	154	305682	44.26 ng/ul 40.13 ng/ul	98	
44) 2-Chloronaphthalene	13.736	162	241676	40.48 ng/ul	97	
45) 2-Nitroaniline	13.936	65	98019	41.33 ng/ul	98	
47) Dimethylphthalate	14.294	163	332302	41.68 ng/ul	99	
48) 2,6-Dinitrotoluene	14.424	165	71932	43.11 ng/ul	95	
50) Acenaphthylene	14.576	152	398117	39.98 ng/ul	99	
51) 3-Nitroaniline	14.759	138	68861	39.90 ng/ul	95	
52) Acenaphthene	14.917	153	267142	40.80 ng/ul	98	
53) 2,4-Dinitrophenol	14.964	184	41668	45.26 ng/ul	92	
55) 4-Nitrophenol	15.058	109	58023	43.76 ng/ul	94	
56) Dibenzofuran	15.246	168	377541	40.29 ng/ul	98	
57) 2,4-Dinitrotoluene	15.211	165	104232	43.77 ng/ul	96	
58) 2,3,4,6-Tetrachlorophenol	15.469	232	78064	46.56 ng/ul	100	
59) Diethylphthalate	15.646	149	356591	41.78 ng/ul	99	
61) Fluorene	15.898	166	301109	40.60 ng/ul	99	
62) 4-Chlorophenyl-phenyle	15.881	204	159698	41.35 ng/ul	98	
63) 4-Nitroaniline	15.922	138	75740	44.24 ng/ul	96	
66) 4,6-Dinitro-2-methylph	15.975	198	62313	46.41 ng/ul#	98	
67) N-Nitrosodiphenylamine 68) 4-Bromophenyl-phenylether	16.098 16.774	169 248	274904 100941	43.31 ng/ul 44.69 ng/ul	99 96	
69) Hexachlorobenzene	16.897		100941	44.09 ng/ul 44.92 ng/ul	90 97	
70) Atrazine	17.038		115242	42.81 ng/ul	98	
71) Pentachlorophenol	17.244	266	58572	54.94 ng/ul	95	
72) Phenanthrene	17.637	178	525470	43.35 ng/ul	100	
74) Anthracene	17.732	178	519247	42.69 ng/ul	100	
75) 1,2,3,4-Tetrachloroben	13.660	216	130595	42.24 ng/uL	97	
76) Pentachlorobenzene	15.164	250	118521	41.37 ng/uL	99	
77) Carbazole	18.002	167	487717	44.74 ng/ul	98	
78) Di-n-butylphthalate	18.531	149	625736	43.68 ng/ul	100	
80) Fluoranthene	19.641	202	644020	43.59 ng/ul	98	
82) Pyrene	19.999	202	622887	43.14 ng/ul	99	
83) Butylbenzylphthalate	20.863	149	274997	44.29 ng/ul	99	
84) 3,3'-Dichlorobenzidine	21.780	252	192302	41.42 ng/ul	96	
85) Benzo(a)anthracene	21.874	228	577066	43.73 ng/ul	99	
86) Bis(2-ethylhexyl)phtha 87) Chrysene	21.739	149 229	386376	43.36 ng/ul	99 100	
89) Di-n-octyl phthalate	21.944 23.008	228 149	552536 649100	43.83 ng/ul 42.59 ng/ul	100 100	
90) Benzo(b)fluoranthene	23.008	252	582869	43.71 ng/ul	99	
91) Benzo(k)fluoranthene	24.200	252	533814	42.66 ng/ul	99	
93) Benzo(a)pyrene	25.135	252	541799	42.66 ng/ul	98	
94) Indeno(1,2,3-cd)pyrene	29.195	276		> 43.17 ng/ul >	ĨĬ	1361 34
95) Dibenzo(a,h)anthracene	29.259		514449	43.00 ng/ul	98	•••
96) Benzo(g,h,i)perylene		276	508350	42.95 ng/ul	96	
				-		

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