





Data Fi Acq On Dperato Sample Misc	: PB141091BS :		ta\BG1	20121\			Instrument : BNA_G ClientSampleId : SLCS091
ALS Via	ol : 6 Sample Multipli	er: 1	Manual IntegrationsAPPROVED				
Quant M Quant T QLast U	ime: Dec 01 18:49:27 202 lethod : Z:\svoasrv\HPCHE itle : SVOA CALIBRATION lpdate : Wed Nov 24 06:04	M1\BNA_G :50 2021	21.M	Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021			
Respons	e via : Initial Calibrat	ion					
	Compound	R.T.	QIon	Response	Conc Un	its Dev	(Min)
Intern	al Standards						
	,4-Dichlorobenzene-d4	8.188	152	28533	20.000	ng/ul	-0.01
	aphthalene-d8	11.020		123394	20.000		0.00
	cenaphthene-d10	14.822		81168	20.000	-	0.00
	henanthrene-d10	17.572		177933	20.000	-	0.00
79) C	hrysene-d12	21.872	240	156456	20.000	-	0.00
88) P	erylene-d12	25.274	264	155833	20.000	-	0.00
						-	
	Monitoring Compounds						
	,4-Dioxane-d8	3.529	96	5148	6.270		-0.01
	yridine-d5	3.958	84	77848	32.311	-	-0.02
	henol-d5	7.354	99	97542	34.589	-	0.00
	is-(2-Chloroethyl)eth	7.507	67	62330	35.193	-	0.00
	-Chlorophenol-d4	7.724	132	70953	34.940	-	0.00
	-Methylphenol-d8	8.905	113	77104	33.882	-	0.00
	itrobenzene-d5	9.369	128	37687	36.181	-	0.00
	-Nitrophenol-d4	10.098	143	42484	36.157	-	0.00
	,4-Dichlorophenol-d3 -Chloroaniline-d4	10.644 11.156	165 131	72277	36.255	-	0.00
	imethylphthalate-d6	14.217	166	90957 223429	31.181	-	0.00
	cenaphthylene-d8	14.522	160	223429	35.775	-	0.00 0.00
	-Nitrophenol-d4	15.045	143	32751	35.922 32.397	-	0.00
	luorene-d10	15.815	176	199423	35.459	-	0.00
	,6-Dinitro-2-methylph	15.950	200	37751	34.383	-	0.00
	nthracene-d10	17.677	188	305916	35.948		0.00
	/rene-d10	19.951	212	351235	37.102		0.00
	enzo(a)pyrene-d12	25.039	264	308018	37.010		0.00
						0	
-	Compounds					Qva	alue
• •	,4-Dioxane	3.570	88	12160	13.132		94
-	/ridine	3.982	79	79005	31.513	ng/ul	97
	enzaldehyde	7.325	77	61747	34.382	-	95
8) Ph		7.378	94	102062	34.936	-	99
	s(2-Chloroethyl)ether	7.601	93	77344	34.995	-	96
	Chlorophenol	7.754	128	72047	34.816	-	97
	Methylphenol	8.641	108	74686	34.322	-	98
	2'-oxybis(1-Chloropr	8.711 9.023	45 105	112413	35.247	-	98
	Nitroso-di-n-propyla	8.993	70	119239 70423	33.875 34.816		98 98
	Methylphenol	8.970	108	79894	34.336		98
	exachloroethane	9.275	117	29195	33.402	-	96
•	trobenzene	9.416	77	102064	37.369	-	97
-	ophorone	9.933	82	194419	36.639	-	99
	Nitrophenol	10.127	139	42909	35.257		98
	4-Dimethylphenol	10.180	107	90659	36.434		97
27) Bi	s(2-Chloroethoxy)met	10.409	93	108267	36.959 (98
29)2,	4-Dichlorophenol	10.674	162	69067	35.194 i	-	97
	phthalene	11.073	128	236042	35.156 (-	98
	Chloroaniline	11.185	127	92171	31.474 ၊	-	97
	xachlorobutadiene	11.332	225	45597	33.686 1	-	98 19 / 9 / 9 / 9 / 9 / 9 / 9 / 9 / 9 / 9 /
	prolactam	11.943	113	27222m>			
33) 4-	Chloro-3-methylphenol	12.301	107	85797	36.395 r	ng/u1	95

Acq On : 1 Dec 2021 18:07 Dperator : CG/JU : PB141091BS Sample **1isc** : ALS Vial : 6 Sample Multiplier: 1 Quant Time: Dec 01 18:49:27 2021 Juant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M Quant Title : SVOA CALIBRATION 2Last Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120121\

Data File : BG051290.D

Instrument : BNA_G ClientSampleId : SLCS091

Manual IntegrationsAPPROVED

Reviewed By : Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.666	142	159480	34.921 ng/ul	98
37) 1-Methylnaphthalene	12.883	142	161317	34.334 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.024	216	91708	35.989 ng/ul	99
40) Hexachlorocyclopentadiene	12.989	237	17085	16.588 ng/ul#	96
41) 2,4,6-Trichlorophenol	13.271	196	56166	35.124 ng/ul	100
42) 2,4,5-Trichlorophenol	13.353	196	59308	35.417 ng/ul	97
43) 1,1'-Biphenyl	13.659	154	216049	35.637 ng/ul	98
44) 2-Chloronaphthalene	13.711	162	169707	35.191 ng/ul	98
45) 2-Nitroaniline	13.917	65	63653	38.138 ng/ul	94
47) Dimethylphthalate	14.264	163	222588	35.211 ng/ul	99
48) 2,6-Dinitrotoluene	14.405	165	47093	35.465 ng/ul	96
50) Acenaphthylene	14.552	152	273095	35.099 ng/ul	99
51) 3-Nitroaniline	14.740	138	47523	36.206 ng/ul	95
52) Acenaphthene	14.886	153	182846	35.633 ng/ul	98
53) 2,4-Dinitrophenol	14.963	184	18135	24.708 ng/ul	88
55) 4-Nitrophenol	15.057	109	28709	32.737 ng/ul	96
56) Dibenzofuran	15.221	168	256190	34.614 ng/ul	97
57) 2,4-Dinitrotoluene	15.198	165	68884	36.320 ng/ul	97
58) 2,3,4,6-Tetrachlorophenol	15.451	232	43136	32.804 ng/ul	98
59) Diethylphthalate	15.621	149	237579	35.804 ng/ul	99
61) Fluorene	15.874	166	206972	34.911 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.850	204	110454	34.571 ng/ul	97
63) 4-Nitroaniline	15.903	138	50656	39.658 ng/ul	96
66) 4,6-Dinitro-2-methylph	15.962	198	35141	33.186 ng/ul#	99
67) N-Nitrosodiphenylamine	16.073	169	185580	36.432 ng/ul	97
68) 4-Bromophenyl-phenylether	16.749	248	67704	35.502 ng/ul	94
69) Hexachlorobenzene	16.872	284	70290	36.147 ng/ul	98
70) Atrazine	17.013	200	70422	32.895 ng/ul	98 100
71) Pentachlorophenol 72) Phenanthrene	17.231 17.619	266 178	18815 352301	21.836 ng/ul	100 99
74) Anthracene	17.713	178	348044	35.860 ng/ul 35.671 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.629	216	94781	36.519 ng/uL	98 99
76) Pentachlorobenzene	15.145	250	82978	34.313 ng/uL	98
77) Carbazole	17.983	167	319258	37.277 ng/ul	98
78) Di-n-butylphthalate	18.506	149	405739	36.742 ng/ul	99
80) Fluoranthene	19.622	202	434186	37.342 ng/ul	98
82) Pyrene	19.981	202	421269	37.038 ng/ul	99
83) Butylbenzylphthalate	20.844	149	178727	37.798 ng/ul	93
84) 3,3'-Dichlorobenzidine	21.761	252	127250	34.933 ng/ul	98
85) Benzo(a)anthracene	21.855	228	384549	36.238 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.714	149	254994	37.476 ng/ul	99
87) Chrysene	21.925	228	370001	36.295 ng/ul	99
89) Di-n-octyl phthalate	22.977	149	432637	38.322 ng/ul	100
90) Benzo(b)fluoranthene	24.187	252	382828	36.402 ng/ul	99
91) Benzo(k)fluoranthene	24.258	252	366533	37.140 ng/ul	99
93) Benzo(a)pyrene	25.110	252	365390	36.418 ng/ul	97
94) Indeno(1,2,3-cd)pyrene	29.193	276	408446	36.380 ng/ul	97
95) Dibenzo(a,h)anthracene	29.240	278	346007	36.327 ng/ul	96
96) Benzo(g,h,i)perylene	30.409	276	346271	36.657 ng/ul	97

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