

Supervised By : Jagrut
 Prepared By : Jagrut
 02/09/2023

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Pag 1

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020823\
 Data File : BF132358.D
 Acq On : 08 Feb 2023 15:47
 Operator : CG\JU
 Sample : 01481-03MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Feb 09 01:59:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 09 01:56:26 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian
 Giraldo

Supervised By : Jagrut
 Spadhyard
 Upd 09/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.366	196	312571	44.629	ng	99	
46) 1,1'-Biphenyl	9.513	154	1089805	46.731	ng	97	
47) 2-Chloronaphthalene	9.542	162	853753	48.066	ng	99	
48) 2-Nitroaniline	9.636	65	208073	39.389	ng	#	78
49) Acenaphthylene	9.960	152	1291498	44.552	ng	99	
50) Dimethylphthalate	9.813	163	994265	46.989	ng	99	
51) 2,6-Dinitrotoluene	9.877	165	224065	47.414	ng	#	83
52) Acenaphthene	10.136	154	872881	48.096	ng	99	
53) 3-Nitroaniline	10.048	138	165520	28.882	ng	92	
54) 2,4-Dinitrophenol	10.148	184	133905	50.565	ng	99	
55) Dibenzofuran	10.307	168	1172505	45.981	ng	100	
56) 4-Nitrophenol	10.201	139	333329	77.293	ng	92	
57) 2,4-Dinitrotoluene	10.283	165	288104	46.932	ng	88	
58) Fluorene	10.654	166	903145	46.021	ng	100	
59) 2,3,4,6-Tetrachlorophenol	10.419	232	233761	45.783	ng	98	
60) Diethylphthalate	10.513	149	985809	46.832	ng	99	
61) 4-Chlorophenyl-phenyle...	10.636	204	429001	46.062	ng	95	
62) 4-Nitroaniline	10.666	138	194810	35.448	ng	96	
63) Azobenzene	10.801	77	821623	40.072	ng	93	
65) 4,6-Dinitro-2-methylph...	10.689	198	93048	33.596	ng	95	
66) n-Nitrosodiphenylamine	10.760	169	763137	49.941	ng	99	
67) 4-Bromophenyl-phenylether	11.130	248	264700	52.971	ng	95	
68) Hexachlorobenzene	11.201	284	288853	54.082	ng	94	
69) Atrazine	11.283	200	244603	55.651	ng	99	
70) Pentachlorophenol	11.395	266	287777	78.418	ng	99	
71) Phenanthrene	11.624	178	1848256	72.479	ng	99	
72) Anthracene	11.672	178	1385739	53.397	ng	100	
73) Carbazole	11.824	167	1010991	43.447	ng	100	
74) Di-n-butylphthalate	12.148	149	1341994	48.782	ng	99	
75) Fluoranthene	12.824	202	2244486	86.251	ng	96	
77) Benzidine	12.936	184	340057	36.911	ng	98	
78) Pyrene	13.054	202	2139008	100.919	ng	98	
80) Butylbenzylphthalate	13.660	149	471410	50.341	ng	94	
81) Benzo(a)anthracene	14.248	228	1542568	83.694	ng	97	
82) 3,3'-Dichlorobenzidine	14.201	252	177300	30.203	ng	99	
83) Chrysene	14.289	228	1305844	75.664	ng	99	
84) Bis(2-ethylhexyl)phtha...	14.213	149	651337	51.459	ng	99	
85) Di-n-octyl phthalate	14.848	149	1063116	51.495	ng	#	96
87) Indeno(1,2,3-cd)pyrene	17.465	276	907547	56.416	ng	95	
88) Benzo(b)fluoranthene	15.365	252	1407375	91.454	ng	98	
89) Benzo(k)fluoranthene	15.395	252	985675	67.102	ng	98	
90) Benzo(a)pyrene	15.765	252	1095334	76.436	ng	99	
91) Dibenzo(a,h)anthracene	17.483	278	589231	45.559	ng	98	
92) Benzo(g,h,i)perylene	17.965	276	756916	56.649	ng	97	

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

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