

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090622\
 Data File : BF130140.D
 Acq On : 06 Sep 2022 16:31
 Operator : CG\JU
 Sample : PB147437BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147437BS

Manual Integrations
 APPROVED

Reviewed By : Christian
 Giraldo

Quant Time: Sep 07 02:13:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 07 02:05:07 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							09/07/2022
1) 1,4-Dichlorobenzene-d4	6.698	152	241533	20.000	ng	0.00	Supervised By :Jagrut padhyay
21) Naphthalene-d8	7.981	136	1013660	20.000	ng	0.00	
39) Acenaphthene-d10	9.733	164	556931	20.000	ng	0.00	
64) Phenanthrene-d10	11.210	188	961044	20.000	ng	0.00	
76) Chrysene-d12	13.839	240	540935	20.000	ng	# 0.00	
86) Perylene-d12	15.233	264	468122	20.000	ng	0.00	09/07/2022
System Monitoring Compounds							
5) 2-Fluorophenol	5.310	112	1721804	121.020	ng	0.01	
7) Phenol-d6	6.334	99	2215874	124.926	ng	0.00	
23) Nitrobenzene-d5	7.269	82	1342932	85.142	ng	0.00	
42) 2,4,6-Tribromophenol	10.522	330	719875	141.998	ng	0.00	
45) 2-Fluorobiphenyl	9.057	172	2581489	76.850	ng	0.00	
79) Terphenyl-d14	12.798	244	2842019	91.109	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	2.428	88	187591	28.213	ng		95
3) Pyridine	3.134	79	521913	29.272	ng		91
4) n-Nitrosodimethylamine	3.075	42	263211	37.539	ng		84
6) Aniline	6.363	93	785040	35.302	ng	#	50
8) 2-Chlorophenol	6.481	128	692625	44.354	ng		99
9) Benzaldehyde	6.245	77	326380	30.063	ng		97
10) Phenol	6.351	94	842737	42.033	ng	#	69
11) bis(2-Chloroethyl)ether	6.439	93	626847	41.150	ng		99
12) 1,3-Dichlorobenzene	6.640	146	667816	38.578	ng		96
13) 1,4-Dichlorobenzene	6.716	146	674834	38.700	ng		96
14) 1,2-Dichlorobenzene	6.863	146	648005	40.881	ng		98
15) Benzyl Alcohol	6.845	79	516801	43.227	ng		93
16) 2,2'-oxybis(1-Chloropr...	6.981	45	809473	39.099	ng	#	52
17) 2-Methylphenol	6.957	107	561160	42.712	ng	#	87
18) Hexachloroethane	7.210	117	247226	39.263	ng		94
19) n-Nitroso-di-n-propyla...	7.122	70	441375	41.305	ng		93
20) 3+4-Methylphenols	7.110	107	651294	42.273	ng	#	75
22) Acetophenone	7.116	105	876293	38.694	ng	#	98
24) Nitrobenzene	7.287	77	675879	40.292	ng		95
25) Isophorone	7.528	82	1325795	41.529	ng		98
26) 2-Nitrophenol	7.598	139	337122	43.162	ng		94
27) 2,4-Dimethylphenol	7.639	122	626703	46.785	ng		98
28) bis(2-Chloroethoxy)met...	7.739	93	813704	42.432	ng		98
29) 2,4-Dichlorophenol	7.839	162	596149	41.606	ng		99
30) 1,2,4-Trichlorobenzene	7.922	180	562748	37.248	ng		98
31) Naphthalene	8.004	128	1915702	37.471	ng		99
32) Benzoic acid	7.775	122	487806	45.066	ng		100
33) 4-Chloroaniline	8.051	127	662112	32.056	ng		98
34) Hexachlorobutadiene	8.122	225	318304	36.787	ng		99
35) Caprolactam	8.434	113	190700m	42.949	ng		
36) 4-Chloro-3-methylphenol	8.539	107	645613	44.528	ng		94
37) 2-Methylnaphthalene	8.692	142	1308140	39.722	ng		98
38) 1-Methylnaphthalene	8.792	142	1244483	38.878	ng		99
40) 1,2,4,5-Tetrachloroben...	8.857	216	550569	37.911	ng		98
41) Hexachlorocyclopentadiene	8.845	237	631074	83.149	ng		96
43) 2,4,6-Trichlorophenol	8.969	196	427700	41.259	ng		98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.010	196	468560	41.703	ng	95	09/07/2022
46) 1,1'-Biphenyl	9.157	154	1578117	39.003	ng	99	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	9.181	162	1244951	38.402	ng	99	
48) 2-Nitroaniline	9.281	65	389542	48.760	ng	91	
49) Acenaphthylene	9.592	152	1941435	39.437	ng	100	
50) Dimethylphthalate	9.469	163	1532620	40.403	ng	100	
51) 2,6-Dinitrotoluene	9.528	165	351032	47.463	ng	# 81	09/07/2022
52) Acenaphthene	9.769	154	1358766m	43.795	ng		
53) 3-Nitroaniline	9.692	138	319282	37.611	ng	99	
54) 2,4-Dinitrophenol	9.798	184	333494	97.367	ng	90	
55) Dibenzofuran	9.939	168	1767069	39.775	ng	99	
56) 4-Nitrophenol	9.857	139	631436	96.481	ng	93	
57) 2,4-Dinitrotoluene	9.928	165	443221	52.051	ng	# 86	
58) Fluorene	10.280	166	1313549	38.858	ng	99	
59) 2,3,4,6-Tetrachlorophenol	10.051	232	368380	42.213	ng	# 78	
60) Diethylphthalate	10.163	149	1509123	41.123	ng	99	
61) 4-Chlorophenyl-phenyle...	10.275	204	593629	39.726	ng	98	
62) 4-Nitroaniline	10.310	138	397377	48.132	ng	98	
63) Azobenzene	10.433	77	1424042	39.412	ng	95	
65) 4,6-Dinitro-2-methylph...	10.327	198	214105	47.408	ng	99	
66) n-Nitrosodiphenylamine	10.398	169	1273870	39.651	ng	98	
67) 4-Bromophenyl-phenylether	10.763	248	428417	40.077	ng	95	
68) Hexachlorobenzene	10.822	284	477928	41.032	ng	92	
69) Atrazine	10.927	200	421166	46.339	ng	98	
70) Pentachlorophenol	11.016	266	558955	85.817	ng	96	
71) Phenanthrene	11.239	178	2018335	38.726	ng	99	
72) Anthracene	11.292	178	2066652	40.337	ng	99	
73) Carbazole	11.445	167	1976466	41.833	ng	99	
74) Di-n-butylphthalate	11.780	149	2244244	39.392	ng	100	
75) Fluoranthene	12.421	202	2082642	40.297	ng	98	
77) Benzidine	12.545	184	883182	60.472	ng	99	
78) Pyrene	12.645	202	2097631	43.233	ng	98	
80) Butylbenzylphthalate	13.274	149	807204	41.827	ng	99	
81) Benzo(a)anthracene	13.827	228	1495829	40.351	ng	99	
82) 3,3'-Dichlorobenzidine	13.798	252	435220	37.647	ng	99	
83) Chrysene	13.868	228	1454318	39.792	ng	100	
84) Bis(2-ethylhexyl)phtha...	13.833	149	808907	33.721	ng	99	
85) Di-n-octyl phthalate	14.445	149	1297432	34.240	ng	98	
87) Indeno(1,2,3-cd)pyrene	16.580	276	1433349	46.418	ng	97	
88) Benzo(b)fluoranthene	14.845	252	1378042	44.143	ng	99	
89) Benzo(k)fluoranthene	14.868	252	1229960	40.738	ng	97	
90) Benzo(a)pyrene	15.180	252	1153749	46.511	ng	98	
91) Dibenzo(a,h)anthracene	16.604	278	1246639	49.338	ng	97	
92) Benzo(g,h,i)perylene	16.992	276	1264389	50.015	ng	# 92	

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

