

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF042622\  
 Data File : BF127907.D  
 Acq On : 26 Apr 2022 11:47  
 Operator : CG\JU  
 Sample : PB144358BSD  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB144358BSD

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 04/27/2022  
 Supervised By : Jagrut Upadhyay 04/27/2022

Quant Time: Apr 27 01:46:25 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF042222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Apr 24 00:43:03 2022  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
Internal Standards							
1) 1,4-Dichlorobenzene-d4	6.945	152	185743	20.000	ng	0.00	
21) Naphthalene-d8	8.228	136	724155	20.000	ng	0.00	
39) Acenaphthene-d10	9.992	164	395301	20.000	ng	0.00	
64) Phenanthrene-d10	11.486	188	661501	20.000	ng	0.01	
76) Chrysene-d12	14.133	240	400561	20.000	ng	0.00	
86) Perylene-d12	15.645	264	345403	20.000	ng	0.01	
System Monitoring Compounds							
5) 2-Fluorophenol	5.587	112	1009712	86.585	ng	0.02	
7) Phenol-d6	6.575	99	1323488	87.465	ng	0.00	
23) Nitrobenzene-d5	7.516	82	1360808	88.366	ng	0.00	
42) 2,4,6-Tribromophenol	10.786	330	348904	87.678	ng	0.00	
45) 2-Fluorobiphenyl	9.310	172	2077860	88.925	ng	0.00	
79) Terphenyl-d14	13.074	244	2029461	92.645	ng	0.01	
Target Compounds							
2) 1,4-Dioxane	2.951	88	196964	35.400	ng		98
3) Pyridine	3.687	79	541917	38.013	ng		98
4) n-Nitrosodimethylamine	3.663	42	324242	46.987	ng		100
6) Aniline	6.610	93	681104	37.885	ng		99
8) 2-Chlorophenol	6.734	128	587073	48.504	ng		99
9) Benzaldehyde	6.498	77	391278	41.799	ng		95
10) Phenol	6.592	94	712415m	46.681	ng		
11) bis(2-Chloroethyl)ether	6.681	93	596127	47.088	ng		98
12) 1,3-Dichlorobenzene	6.886	146	613181	44.012	ng		97
13) 1,4-Dichlorobenzene	6.963	146	612441	44.040	ng		97
14) 1,2-Dichlorobenzene	7.116	146	591534	45.948	ng		99
15) Benzyl Alcohol	7.092	79	489843	47.604	ng		98
16) 2,2'-oxybis(1-Chloropr...	7.216	45	904707	46.757	ng		97
17) 2-Methylphenol	7.198	107	534506	51.183	ng		96
18) Hexachloroethane	7.457	117	243178	47.565	ng		99
19) n-Nitroso-di-n-propyla...	7.363	70	441591	47.674	ng		99
20) 3+4-Methylphenols	7.351	107	551893	43.748	ng		90
22) Acetophenone	7.357	105	783074	44.367	ng	#	91
24) Nitrobenzene	7.533	77	687183	46.309	ng		98
25) Isophorone	7.769	82	1284011	49.540	ng		100
26) 2-Nitrophenol	7.845	139	316283	50.018	ng		97
27) 2,4-Dimethylphenol	7.880	122	548413	52.215	ng		98
28) bis(2-Chloroethoxy)met...	7.975	93	756701	45.555	ng		100
29) 2,4-Dichlorophenol	8.086	162	494380	44.930	ng		99
30) 1,2,4-Trichlorobenzene	8.169	180	550097	44.012	ng		98
31) Naphthalene	8.251	128	1638867	44.435	ng		99
32) Benzoic acid	8.016	122	299606	39.798	ng		98
33) 4-Chloroaniline	8.298	127	348830	23.904	ng		99
34) Hexachlorobutadiene	8.363	225	352439	44.767	ng		98
35) Caprolactam	8.686	113	144147m	40.584	ng		
36) 4-Chloro-3-methylphenol	8.786	107	533842	45.428	ng		99
37) 2-Methylnaphthalene	8.945	142	1092075	44.803	ng		99
38) 1-Methylnaphthalene	9.045	142	1038262	43.823	ng		99
40) 1,2,4,5-Tetrachloroben...	9.110	216	508838	43.671	ng		98
41) Hexachlorocyclopentadiene	9.092	237	706250	111.877	ng		99
43) 2,4,6-Trichlorophenol	9.222	196	349842	45.632	ng		100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.263	196	361181	43.337	ng	99
46) 1,1'-Biphenyl	9.410	154	1312132	44.811	ng	99
47) 2-Chloronaphthalene	9.439	162	966950	43.378	ng	99
48) 2-Nitroaniline	9.533	65	352005	46.683	ng	94
49) Acenaphthylene	9.857	152	1542091	45.225	ng	100
50) Dimethylphthalate	9.710	163	1151390	43.431	ng	99
51) 2,6-Dinitrotoluene	9.774	165	265743	46.297	ng	95
52) Acenaphthene	10.027	154	1102252m	48.082	ng	
53) 3-Nitroaniline	9.945	138	167085	26.214	ng	98
54) 2,4-Dinitrophenol	10.057	184	276311	91.406	ng	# 83
55) Dibenzofuran	10.198	168	1365349	41.281	ng	98
56) 4-Nitrophenol	10.116	139	374374	76.918	ng	96
57) 2,4-Dinitrotoluene	10.186	165	346665	45.595	ng	96
58) Fluorene	10.545	166	1033741	41.886	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.316	232	300652	42.899	ng	99
60) Diethylphthalate	10.410	149	1115375	42.905	ng	99
61) 4-Chlorophenyl-phenyle...	10.533	204	517535	42.199	ng	97
62) 4-Nitroaniline	10.569	138	256255	41.123	ng	98
63) Azobenzene	10.692	77	1193913	41.737	ng	100
65) 4,6-Dinitro-2-methylph...	10.592	198	183829	46.769	ng	97
66) n-Nitrosodiphenylamine	10.651	169	925886	44.983	ng	99
67) 4-Bromophenyl-phenylether	11.021	248	331669	45.229	ng	99
68) Hexachlorobenzene	11.092	284	359339	46.510	ng	# 91
69) Atrazine	11.180	200	328378	51.593	ng	98
70) Pentachlorophenol	11.286	266	385467	84.806	ng	97
71) Phenanthrene	11.510	178	1458620	42.124	ng	99
72) Anthracene	11.563	178	1503538	43.590	ng	99
73) Carbazole	11.716	167	1341015	40.339	ng	99
74) Di-n-butylphthalate	12.033	149	1636035	43.766	ng	99
75) Fluoranthene	12.704	202	1516038	41.600	ng	98
77) Benzidine	12.821	184	703471	92.816	ng	98
78) Pyrene	12.933	202	1519875	47.004	ng	99
80) Butylbenzylphthalate	13.539	149	637264	50.281	ng	100
81) Benzo(a)anthracene	14.121	228	1229737	46.060	ng	99
82) 3,3'-Dichlorobenzidine	14.080	252	254483	30.020	ng	100
83) Chrysene	14.162	228	1108636	43.479	ng	99
84) Bis(2-ethylhexyl)phtha...	14.092	149	861539	51.260	ng	100
85) Di-n-octyl phthalate	14.715	149	1351453	51.887	ng	100
87) Indeno(1,2,3-cd)pyrene	17.203	276	1036751	44.969	ng	99
88) Benzo(b)fluoranthene	15.198	252	1093628	50.411	ng	99
89) Benzo(k)fluoranthene	15.233	252	915135	41.870	ng	99
90) Benzo(a)pyrene	15.586	252	966955	53.734	ng	99
91) Dibenzo(a,h)anthracene	17.227	278	892029	48.244	ng	98
92) Benzo(g,h,i)perylene	17.686	276	916781	47.049	ng	99

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

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