

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122422\
 Data File : BF131826.D
 Acq On : 24 Dec 2022 15:35
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/28/2022
 Supervised By : Yogesh Patel 12/28/2022

Quant Time: Dec 26 04:38:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 26 04:31:10 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	76108	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	251276	20.000	ng	0.00
39) Acenaphthene-d10	9.892	164	151284	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	296782	20.000	ng	0.00
76) Chrysene-d12	14.045	240	216611	20.000	ng	# 0.00
86) Perylene-d12	15.521	264	143252	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	540020	118.372	ng	0.00
7) Phenol-d6	6.475	99	709600	118.697	ng	0.00
23) Nitrobenzene-d5	7.416	82	767063	115.412	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	202506	120.481	ng	0.00
45) 2-Fluorobiphenyl	9.216	172	1280779	106.746	ng	0.00
79) Terphenyl-d14	12.980	244	1609045	101.872	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.528	88	124443	60.433	ng	98
3) Pyridine	3.281	79	364697	63.791	ng	97
4) n-Nitrosodimethylamine	3.263	42	170456	54.062	ng	98
6) Aniline	6.504	93	429570	60.082	ng	98
8) 2-Chlorophenol	6.622	128	285233	58.998	ng	94
9) Benzaldehyde	6.387	77	204384	49.906	ng	99
10) Phenol	6.487	94	421718	63.242	ng	91
11) bis(2-Chloroethyl)ether	6.575	93	303839	60.785	ng	98
12) 1,3-Dichlorobenzene	6.775	146	320361	57.154	ng	98
13) 1,4-Dichlorobenzene	6.857	146	329875	58.346	ng	96
14) 1,2-Dichlorobenzene	7.010	146	301367	55.800	ng	98
15) Benzyl Alcohol	6.987	79	310961	56.737	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.116	45	379421	58.762	ng	99
17) 2-Methylphenol	7.098	107	262642	60.279	ng	98
18) Hexachloroethane	7.351	117	131674	57.734	ng	90
19) n-Nitroso-di-n-propyla...	7.263	70	246763	55.048	ng	99
20) 3+4-Methylphenols	7.257	107	334444	56.355	ng	96
22) Acetophenone	7.257	105	446368	56.686	ng	# 98
24) Nitrobenzene	7.434	77	385915	57.570	ng	98
25) Isophorone	7.675	82	633862	59.209	ng	100
26) 2-Nitrophenol	7.745	139	153825	67.291	ng	97
27) 2,4-Dimethylphenol	7.786	122	214922	61.339	ng	98
28) bis(2-Chloroethoxy)met...	7.881	93	333963	60.303	ng	99
29) 2,4-Dichlorophenol	7.992	162	260523	60.134	ng	97
30) 1,2,4-Trichlorobenzene	8.069	180	301837	56.811	ng	100
31) Naphthalene	8.151	128	828005	58.505	ng	99
32) Benzoic acid	7.934	122	180943m	71.307	ng	
33) 4-Chloroaniline	8.204	127	322392	60.772	ng	97
34) Hexachlorobutadiene	8.263	225	206105	55.043	ng	99
35) Caprolactam	8.604	113	78839m	70.348	ng	
36) 4-Chloro-3-methylphenol	8.692	107	291654	59.951	ng	98
37) 2-Methylnaphthalene	8.845	142	555203	57.363	ng	99
38) 1-Methylnaphthalene	8.945	142	545418	58.702	ng	100
40) 1,2,4,5-Tetrachloroben...	9.010	216	311506	55.624	ng	100
41) Hexachlorocyclopentadiene	8.992	237	137474	48.364	ng	99
43) 2,4,6-Trichlorophenol	9.128	196	202269	58.393	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.169	196	218507	58.656	ng	97
46) 1,1'-Biphenyl	9.316	154	674064	56.354	ng	98
47) 2-Chloronaphthalene	9.339	162	558960	57.356	ng	99
48) 2-Nitroaniline	9.439	65	214407	62.107	ng	93
49) Acenaphthylene	9.757	152	833074	58.220	ng	100
50) Dimethylphthalate	9.622	163	675818	58.600	ng	100
51) 2,6-Dinitrotoluene	9.686	165	147141	61.801	ng	99
52) Acenaphthene	9.928	154	512274	53.440	ng	98
53) 3-Nitroaniline	9.863	138	155321	67.994	ng	88
54) 2,4-Dinitrophenol	9.963	184	93585	78.307	ng	94
55) Dibenzofuran	10.104	168	773188	57.086	ng	100
56) 4-Nitrophenol	10.022	139	111595	70.322	ng	91
57) 2,4-Dinitrotoluene	10.092	165	205790	65.159	ng	98
58) Fluorene	10.445	166	637899	57.933	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.222	232	181670m	57.590	ng	
60) Diethylphthalate	10.322	149	649157	58.656	ng	100
61) 4-Chlorophenyl-phenyle...	10.439	204	334632	54.994	ng	93
62) 4-Nitroaniline	10.480	138	158697	70.106	ng	97
63) Azobenzene	10.598	77	718563	56.926	ng	97
65) 4,6-Dinitro-2-methylph...	10.504	198	130757	69.561	ng	92
66) n-Nitrosodiphenylamine	10.563	169	540236	56.306	ng	99
67) 4-Bromophenyl-phenylether	10.927	248	205837	53.914	ng	94
68) Hexachlorobenzene	10.992	284	224791	54.433	ng	95
69) Atrazine	11.092	200	167427	54.951	ng	99
70) Pentachlorophenol	11.186	266	121952	57.362	ng	99
71) Phenanthrene	11.416	178	962874	58.207	ng	100
72) Anthracene	11.469	178	950650	59.068	ng	100
73) Carbazole	11.621	167	847578	64.520	ng	99
74) Di-n-butylphthalate	11.951	149	1022209	63.427	ng	99
75) Fluoranthene	12.610	202	1123629	65.963	ng	99
77) Benzidine	12.727	184	103224	17.785	ng	98
78) Pyrene	12.839	202	1146715	53.688	ng	100
80) Butylbenzylphthalate	13.457	149	427913	69.228	ng	97
81) Benzo(a)anthracene	14.033	228	937609	59.939	ng	100
82) 3,3'-Dichlorobenzidine	13.998	252	253550	56.970	ng	96
83) Chrysene	14.074	228	875546	58.874	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	498673	66.629	ng	99
85) Di-n-octyl phthalate	14.633	149	644083	51.088	ng	100
87) Indeno(1,2,3-cd)pyrene	17.033	276	658919	56.362	ng	100
88) Benzo(b)fluoranthene	15.092	252	615521	65.193	ng	100
89) Benzo(k)fluoranthene	15.121	252	630914	64.022	ng	99
90) Benzo(a)pyrene	15.462	252	493926	61.301	ng	98
91) Dibenzo(a,h)anthracene	17.051	278	538302	55.946	ng	99
92) Benzo(g,h,i)perylene	17.492	276	555082	57.401	ng	97

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

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