

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF043024\
 Data File : BF137771.D
 Acq On : 30 Apr 2024 14:45
 Operator : CG\JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024
 Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 02:59:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 02:54:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.063	152	232068	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	899995	20.000	ng	0.00
39) Acenaphthene-d10	10.110	164	504470	20.000	ng	0.00
64) Phenanthrene-d10	11.604	188	888980	20.000	ng	0.00
76) Chrysene-d12	14.257	240	559359	20.000	ng	0.00
86) Perylene-d12	15.821	264	431596	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	2023763	146.007	ng	0.00
7) Phenol-d6	6.687	99	2554000	145.271	ng	0.01
23) Nitrobenzene-d5	7.634	82	2359121	152.290	ng	0.01
42) 2,4,6-Tribromophenol	10.904	330	813652	157.967	ng	0.00
45) 2-Fluorobiphenyl	9.422	172	4146983	133.366	ng	0.00
79) Terphenyl-d14	13.198	244	4921290	147.444	ng	0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.963	88	486433	77.934	ng	100
3) Pyridine	3.740	79	1172662	77.628	ng	99
4) n-Nitrosodimethylamine	3.716	42	534120	78.520	ng	98
6) Aniline	6.769	93	1281365m	73.253	ng	
8) 2-Chlorophenol	6.851	128	1108715	73.399	ng	99
10) Phenol	6.698	94	1366097	75.833	ng	99
11) bis(2-Chloroethyl)ether	6.798	93	1077065	76.357	ng	98
12) 1,3-Dichlorobenzene	7.004	146	1249151	73.334	ng	99
13) 1,4-Dichlorobenzene	7.081	146	1252452	73.278	ng	99
14) 1,2-Dichlorobenzene	7.239	146	1141543	70.912	ng	100
15) Benzyl Alcohol	7.204	79	936530	77.030	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.328	45	1480936	74.694	ng	100
17) 2-Methylphenol	7.298	107	935639	76.764	ng	98
18) Hexachloroethane	7.575	117	439126	75.999	ng	95
19) n-Nitroso-di-n-propyla...	7.481	70	757565	72.104	ng	96
20) 3+4-Methylphenols	7.457	107	1155345	72.048	ng	93
22) Acetophenone	7.475	105	1470984	72.673	ng	98
24) Nitrobenzene	7.651	77	1213956	76.035	ng	95
25) Isophorone	7.886	82	2183579	78.180	ng	98
26) 2-Nitrophenol	7.957	139	625573	86.525	ng	91
27) 2,4-Dimethylphenol	7.986	122	804579	77.299	ng	98
28) bis(2-Chloroethoxy)met...	8.086	93	1270397	75.123	ng	99
29) 2,4-Dichlorophenol	8.198	162	976304	77.602	ng	99
30) 1,2,4-Trichlorobenzene	8.286	180	1031971	75.244	ng	100
31) Naphthalene	8.369	128	3192837	71.387	ng	98
32) Benzoic acid	8.128	122	826650	92.059	ng	97
33) 4-Chloroaniline	8.428	127	1289903	77.773	ng	99
34) Hexachlorobutadiene	8.481	225	636361	75.574	ng	99
35) Caprolactam	8.804	113	326533m	81.916	ng	
36) 4-Chloro-3-methylphenol	8.886	107	1021112	77.830	ng	95
37) 2-Methylnaphthalene	9.063	142	2079503	71.827	ng	100
38) 1-Methylnaphthalene	9.163	142	2025019	71.096	ng	99
40) 1,2,4,5-Tetrachloroben...	9.222	216	1003395	75.098	ng	100
41) Hexachlorocyclopentadiene	9.210	237	457601	84.319	ng	100
43) 2,4,6-Trichlorophenol	9.333	196	721562	79.336	ng	100
44) 2,4,5-Trichlorophenol	9.375	196	781229	77.791	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.528	154	2486262	71.814	ng	99
47) 2-Chloronaphthalene	9.557	162	2054254	73.495	ng	99
48) 2-Nitroaniline	9.645	65	634046	83.641	ng	98
49) Acenaphthylene	9.975	152	2952810	73.029	ng	98
50) Dimethylphthalate	9.828	163	2419751	75.981	ng	100
51) 2,6-Dinitrotoluene	9.892	165	587445	80.715	ng	89
52) Acenaphthene	10.145	154	1995793	74.467	ng	100
53) 3-Nitroaniline	10.063	138	624809	81.980	ng	95
54) 2,4-Dinitrophenol	10.163	184	338641	94.696	ng	93
55) Dibenzofuran	10.316	168	2809981	72.048	ng	99
56) 4-Nitrophenol	10.210	139	520146	82.343	ng	97
57) 2,4-Dinitrotoluene	10.298	165	797608	83.887	ng	93
58) Fluorene	10.663	166	2203121	70.977	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.427	232	636317	78.045	ng	99
60) Diethylphthalate	10.527	149	2253241	74.496	ng	98
61) 4-Chlorophenyl-phenyle...	10.651	204	1111441	72.671	ng	96
62) 4-Nitroaniline	10.686	138	616892	79.799	ng	96
63) Azobenzene	10.810	77	2154065	73.686	ng	98
65) 4,6-Dinitro-2-methylph...	10.704	198	461569	90.460	ng	98
66) n-Nitrosodiphenylamine	10.769	169	1960910	73.689	ng	99
67) 4-Bromophenyl-phenylether	11.145	248	720367	75.716	ng	94
68) Hexachlorobenzene	11.210	284	787525	75.746	ng	94
69) Atrazine	11.292	200	595932	73.446	ng	98
70) Pentachlorophenol	11.398	266	584621	81.903	ng	99
71) Phenanthrene	11.633	178	3086299	71.227	ng	98
72) Anthracene	11.686	178	3079683	71.432	ng	98
73) Carbazole	11.833	167	3003335	72.282	ng	98
74) Di-n-butylphthalate	12.151	149	3338837	73.664	ng	99
75) Fluoranthene	12.827	202	3407776	71.947	ng	99
77) Benzidine	12.939	184	982924	75.980	ng	98
78) Pyrene	13.057	202	3501594	78.953	ng	98
80) Butylbenzylphthalate	13.663	149	1256487	89.292	ng	95
81) Benzo(a)anthracene	14.245	228	2771657	77.472	ng	99
82) 3,3'-Dichlorobenzidine	14.204	252	747192	69.764	ng	99
83) Chrysene	14.286	228	2580443	77.032	ng	99
84) Bis(2-ethylhexyl)phtha...	14.215	149	1533022	86.444	ng #	97
85) Di-n-octyl phthalate	14.845	149	2098779	87.937	ng	100
87) Indeno(1,2,3-cd)pyrene	17.474	276	2295943	93.569	ng	99
88) Benzo(b)fluoranthene	15.357	252	2101765	78.270	ng	100
89) Benzo(k)fluoranthene	15.392	252	2000880	77.093	ng	99
90) Benzo(a)pyrene	15.762	252	1757069	80.205	ng	99
91) Dibenzo(a,h)anthracene	17.498	278	1828189	91.534	ng	98
92) Benzo(g,h,i)perylene	17.986	276	2020903	96.519	ng	99

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

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