

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011023\
 Data File : BF131988.D
 Acq On : 10 Jan 2023 09:56
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/11/2023
 Supervised By : Jagrut Upadhyay 01/11/2023

Quant Time: Jan 11 00:33:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 10 00:32:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	94676	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	333805	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	195711	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	378485	20.000	ng	0.00
76) Chrysene-d12	13.986	240	223633	20.000	ng	0.00
86) Perylene-d12	15.445	264	163140	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	448101	77.989	ng	0.00
7) Phenol-d6	6.434	99	592908	77.829	ng	0.00
23) Nitrobenzene-d5	7.369	82	646854	80.426	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	184527	82.003	ng	0.00
45) 2-Fluorobiphenyl	9.163	172	1163903	80.283	ng	0.00
79) Terphenyl-d14	12.927	244	1311471	84.053	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.469	88	95709	38.140	ng	99
3) Pyridine	3.199	79	279593	39.699	ng	99
4) n-Nitrosodimethylamine	3.181	42	134420	40.611	ng	91
6) Aniline	6.457	93	389429	38.852	ng	92
8) 2-Chlorophenol	6.575	128	231256	40.116	ng	100
9) Benzaldehyde	6.339	77	189611	37.406	ng	99
10) Phenol	6.451	94	332519	39.128	ng	92
11) bis(2-Chloroethyl)ether	6.534	93	254354	39.704	ng	99
12) 1,3-Dichlorobenzene	6.728	146	251806	39.217	ng	99
13) 1,4-Dichlorobenzene	6.804	146	257040	39.565	ng	98
14) 1,2-Dichlorobenzene	6.963	146	241998	39.221	ng	99
15) Benzyl Alcohol	6.945	79	264545	39.878	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.069	45	314167	38.860	ng	88
17) 2-Methylphenol	7.057	107	224849	39.414	ng	96
18) Hexachloroethane	7.298	117	107742	40.102	ng	100
19) n-Nitroso-di-n-propyla...	7.216	70	214428	38.911	ng	99
20) 3+4-Methylphenols	7.216	107	285507	38.982	ng	90
22) Acetophenone	7.210	105	385122	39.626	ng	99
24) Nitrobenzene	7.386	77	322310	39.364	ng	98
25) Isophorone	7.628	82	602046	40.069	ng	99
26) 2-Nitrophenol	7.698	139	132561	41.146	ng	95
27) 2,4-Dimethylphenol	7.739	122	217885	40.517	ng	99
28) bis(2-Chloroethoxy)met...	7.833	93	329842	39.665	ng	98
29) 2,4-Dichlorophenol	7.945	162	220481	40.028	ng	96
30) 1,2,4-Trichlorobenzene	8.022	180	253953	40.257	ng	98
31) Naphthalene	8.104	128	707657	40.115	ng	99
32) Benzoic acid	7.898	122	124273m	39.861	ng	
33) 4-Chloroaniline	8.157	127	298902	41.060	ng	98
34) Hexachlorobutadiene	8.216	225	177638	40.615	ng	99
35) Caprolactam	8.557	113	63369m	39.171	ng	
36) 4-Chloro-3-methylphenol	8.651	107	255163	39.469	ng	98
37) 2-Methylnaphthalene	8.798	142	510415	40.286	ng	98
38) 1-Methylnaphthalene	8.898	142	472886	40.286	ng	98
40) 1,2,4,5-Tetrachloroben...	8.963	216	285052	41.333	ng	99
41) Hexachlorocyclopentadiene	8.945	237	107392	36.053	ng	100
43) 2,4,6-Trichlorophenol	9.080	196	171839	41.396	ng	99

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44) 2,4,5-Trichlorophenol	9.128	196	195028	40.356	ng	96
46) 1,1'-Biphenyl	9.263	154	613966	40.645	ng	98
47) 2-Chloronaphthalene	9.292	162	474065	40.435	ng	98
48) 2-Nitroaniline	9.392	65	175257	41.156	ng	98
49) Acenaphthylene	9.704	152	734862	40.391	ng	100
50) Dimethylphthalate	9.575	163	614898	40.307	ng	99
51) 2,6-Dinitrotoluene	9.639	165	130621	40.710	ng	94
52) Acenaphthene	9.880	154	438766	40.176	ng	97
53) 3-Nitroaniline	9.810	138	129805	41.699	ng	95
54) 2,4-Dinitrophenol	9.922	184	66667	41.382	ng	90
55) Dibenzofuran	10.051	168	690939	40.137	ng	99
56) 4-Nitrophenol	9.980	139	69695	39.749	ng	99
57) 2,4-Dinitrotoluene	10.045	165	171639	40.011	ng	97
58) Fluorene	10.392	166	549386	39.951	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.169	232	148737	41.305	ng	92
60) Diethylphthalate	10.269	149	611148	40.741	ng	99
61) 4-Chlorophenyl-phenyle...	10.386	204	316328	40.183	ng	98
62) 4-Nitroaniline	10.433	138	116183	41.508	ng	97
63) Azobenzene	10.545	77	643710	40.452	ng	99
65) 4,6-Dinitro-2-methylph...	10.457	198	101855	41.731	ng	88
66) n-Nitrosodiphenylamine	10.510	169	478352	40.285	ng	99
67) 4-Bromophenyl-phenylether	10.874	248	191714	40.491	ng	99
68) Hexachlorobenzene	10.939	284	189572	39.580	ng	95
69) Atrazine	11.039	200	175012	40.530	ng	99
70) Pentachlorophenol	11.139	266	84965	35.811	ng	95
71) Phenanthrene	11.357	178	789661	40.022	ng	99
72) Anthracene	11.410	178	815407	40.144	ng	99
73) Carbazole	11.569	167	670630	39.880	ng	99
74) Di-n-butylphthalate	11.898	149	948693	40.517	ng	99
75) Fluoranthene	12.551	202	873208	39.352	ng	100
77) Benzidine	12.674	184	313405	42.183	ng	99
78) Pyrene	12.780	202	868658	42.588	ng	100
80) Butylbenzylphthalate	13.398	149	349561	42.459	ng	99
81) Benzo(a)anthracene	13.974	228	651359	40.034	ng	98
82) 3,3'-Dichlorobenzidine	13.939	252	207669	40.897	ng	100
83) Chrysene	14.015	228	607480	39.952	ng	99
84) Bis(2-ethylhexyl)phtha...	13.957	149	469201	42.066	ng	99
85) Di-n-octyl phthalate	14.574	149	625029	40.950	ng	98
87) Indeno(1,2,3-cd)pyrene	16.915	276	429542	42.931	ng	100
88) Benzo(b)fluoranthene	15.021	252	471620	42.149	ng	99
89) Benzo(k)fluoranthene	15.051	252	418468	37.473	ng	99
90) Benzo(a)pyrene	15.386	252	404651	40.138	ng	99
91) Dibenzo(a,h)anthracene	16.927	278	360064	42.930	ng	99
92) Benzo(g,h,i)perylene	17.356	276	347339	39.360	ng	97

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

