

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112122\
 Data File : BF131319.D
 Acq On : 21 Nov 2022 17:42
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/22/2022
 Supervised By : Jagrut Upadhyay 11/22/2022

Quant Time: Nov 22 05:33:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 01:17:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	133681	20.000	ng	0.00
21) Naphthalene-d8	8.263	136	472856	20.000	ng	# 0.00
39) Acenaphthene-d10	10.027	164	265524	20.000	ng	0.00
64) Phenanthrene-d10	11.521	188	451716	20.000	ng	0.00
76) Chrysene-d12	14.180	240	272341	20.000	ng	0.00
86) Perylene-d12	15.721	264	226179	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.569	112	538458	76.660	ng	0.00
7) Phenol-d6	6.586	99	678867	75.752	ng	0.00
23) Nitrobenzene-d5	7.539	82	713220	76.032	ng	0.00
42) 2,4,6-Tribromophenol	10.822	330	182629	81.754	ng	0.00
45) 2-Fluorobiphenyl	9.339	172	1298244	71.131	ng	0.00
79) Terphenyl-d14	13.115	244	1266296	76.034	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.787	88	124931	37.308	ng	99
3) Pyridine	3.551	79	350674	37.714	ng	98
4) n-Nitrosodimethylamine	3.528	42	207960	37.926	ng	99
6) Aniline	6.634	93	434825m	37.022	ng	
8) 2-Chlorophenol	6.751	128	306477	38.514	ng	99
9) Benzaldehyde	6.522	77	216206	39.029	ng	93
10) Phenol	6.598	94	377770	38.020	ng	100
11) bis(2-Chloroethyl)ether	6.704	93	309589	37.189	ng	99
12) 1,3-Dichlorobenzene	6.910	146	347480	36.639	ng	99
13) 1,4-Dichlorobenzene	6.986	146	351878	36.514	ng	99
14) 1,2-Dichlorobenzene	7.139	146	324122	36.528	ng	97
15) Benzyl Alcohol	7.110	79	283008	38.514	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.245	45	581112	36.500	ng	99
17) 2-Methylphenol	7.210	107	262980	38.774	ng	99
18) Hexachloroethane	7.486	117	131916	38.266	ng	98
19) n-Nitroso-di-n-propyla...	7.386	70	240593	36.673	ng	99
20) 3+4-Methylphenols	7.369	107	324639	37.488	ng	97
22) Acetophenone	7.381	105	442702	36.781	ng	99
24) Nitrobenzene	7.557	77	365720	37.420	ng	100
25) Isophorone	7.798	82	628203	37.537	ng	98
26) 2-Nitrophenol	7.875	139	135497	42.403	ng	91
27) 2,4-Dimethylphenol	7.904	122	267293	41.199	ng	98
28) bis(2-Chloroethoxy)met...	8.004	93	352280	37.061	ng	99
29) 2,4-Dichlorophenol	8.110	162	278578	39.585	ng	99
30) 1,2,4-Trichlorobenzene	8.198	180	319602	36.660	ng	99
31) Naphthalene	8.280	128	920156	36.517	ng	100
32) Benzoic acid	8.022	122	148782	39.620	ng	86
33) 4-Chloroaniline	8.328	127	356944	35.906	ng	99
34) Hexachlorobutadiene	8.398	225	214540	38.004	ng	98
35) Caprolactam	8.710	113	76665	41.457	ng	92
36) 4-Chloro-3-methylphenol	8.804	107	287393	39.307	ng	94
37) 2-Methylnaphthalene	8.975	142	620507	36.344	ng	98
38) 1-Methylnaphthalene	9.075	142	597061	36.063	ng	100
40) 1,2,4,5-Tetrachloroben...	9.139	216	315399	36.604	ng	99
41) Hexachlorocyclopentadiene	9.127	237	156029	40.281	ng	99
43) 2,4,6-Trichlorophenol	9.251	196	198900	40.691	ng	99

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44) 2,4,5-Trichlorophenol	9.286	196	217528	39.575	ng	95	11/22/2022
46) 1,1'-Biphenyl	9.445	154	739859	36.513	ng	99	Supervised By : Jagrut Upadhyay
47) 2-Chloronaphthalene	9.469	162	595530	36.638	ng	98	
48) 2-Nitroaniline	9.563	65	202449	41.020	ng	96	
49) Acenaphthylene	9.886	152	895425	36.137	ng	99	
50) Dimethylphthalate	9.745	163	695885	37.615	ng	99	
51) 2,6-Dinitrotoluene	9.810	165	144154	38.963	ng	# 79	11/22/2022
52) Acenaphthene	10.063	154	565411	36.441	ng	99	
53) 3-Nitroaniline	9.980	138	143295	37.692	ng	93	
54) 2,4-Dinitrophenol	10.074	184	54839	42.285	ng	98	
55) Dibenzofuran	10.233	168	797493	35.726	ng	99	
56) 4-Nitrophenol	10.122	139	109387	41.459	ng	98	
57) 2,4-Dinitrotoluene	10.210	165	185743	39.901	ng	98	
58) Fluorene	10.580	166	617305	35.434	ng	98	
59) 2,3,4,6-Tetrachlorophenol	10.345	232	188423	41.163	ng	98	
60) Diethylphthalate	10.445	149	659736	37.768	ng	99	
61) 4-Chlorophenyl-phenyle...	10.569	204	317439	35.544	ng	96	
62) 4-Nitroaniline	10.598	138	141641	37.436	ng	99	
63) Azobenzene	10.727	77	672404	35.799	ng	99	
65) 4,6-Dinitro-2-methylph...	10.622	198	77860	41.933	ng	# 76	
66) n-Nitrosodiphenylamine	10.686	169	540461	37.548	ng	99	
67) 4-Bromophenyl-phenylether	11.063	248	207498	37.479	ng	99	
68) Hexachlorobenzene	11.121	284	219574	37.368	ng	97	
69) Atrazine	11.216	200	175165	39.525	ng	99	
70) Pentachlorophenol	11.316	266	125669m	39.440	ng		
71) Phenanthrene	11.551	178	888641	36.402	ng	99	
72) Anthracene	11.598	178	884312	36.859	ng	99	
73) Carbazole	11.751	167	756125	36.947	ng	99	
74) Di-n-butylphthalate	12.080	149	966261	40.069	ng	99	
75) Fluoranthene	12.745	202	934599	36.332	ng	98	
77) Benzidine	12.863	184	162507	32.304	ng	99	
78) Pyrene	12.974	202	927995	38.621	ng	100	
80) Butylbenzylphthalate	13.592	149	329709	41.409	ng	99	
81) Benzo(a)anthracene	14.168	228	697445	37.361	ng	99	
82) 3,3'-Dichlorobenzidine	14.133	252	194640	39.293	ng	# 97	
83) Chrysene	14.210	228	676055	36.913	ng	99	
84) Bis(2-ethylhexyl)phtha...	14.157	149	421293	39.952	ng	# 99	
85) Di-n-octyl phthalate	14.786	149	612888	41.867	ng	99	
87) Indeno(1,2,3-cd)pyrene	17.303	276	610071	40.160	ng	99	
88) Benzo(b)fluoranthene	15.262	252	553176	36.816	ng	100	
89) Benzo(k)fluoranthene	15.292	252	576041	37.306	ng	99	
90) Benzo(a)pyrene	15.656	252	471469	38.241	ng	98	
91) Dibenzo(a,h)anthracene	17.327	278	504127	40.196	ng	98	
92) Benzo(g,h,i)perylene	17.786	276	499989	39.943	ng	98	

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

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