

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092717\
 Data File : BF098892.D
 Acq On : 28 Sep 2017 6:04
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 9/28/2017 2:32:34 PM

Quant Time: Sep 28 07:54:24 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	135315	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	551973	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	229078	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	366422	20.00	ng	0.00
75) Chrysene-d12	14.09	240	280213	20.00	ng	0.00
86) Perylene-d12	15.57	264	214049	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	687798	80.92	ng	-0.01
7) Phenol-d6	6.54	99	843807	80.47	ng	0.00
23) Nitrobenzene-d5	7.48	82	718515	83.48	ng	0.00
41) 2,4,6-Tribromophenol	10.75	330	135251	72.15	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	1175475	85.66	ng	0.00
78) Terphenyl-d14	13.03	244	965884	78.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	162493	40.018	ng	99
3) Pyridine	3.48	79	449021	40.878	ng	98
4) n-Nitrosodimethylamine	3.43	42	191022	41.010	ng	# 96
6) Aniline	6.58	93	550790	38.918	ng	99
8) 2-Chlorophenol	6.70	128	383059	39.794	ng	95
9) Benzaldehyde	6.47	77	230905	37.961	ng	97
10) Phenol	6.56	94	478204	40.777	ng	97
11) bis(2-Chloroethyl)ether	6.65	93	355406	38.983	ng	98
12) 1,3-Dichlorobenzene	6.86	146	398707	39.549	ng	99
13) 1,4-Dichlorobenzene	6.93	146	406071	39.590	ng	100
14) 1,2-Dichlorobenzene	7.09	146	375622	39.633	ng	100
15) Benzyl Alcohol	7.06	79	302993	39.388	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.19	45	547465	38.542	ng	99
17) 2-Methylphenol	7.16	107	304852	38.705	ng	98
18) Hexachloroethane	7.43	117	119608	32.442	ng	93
19) n-Nitroso-di-n-propylamine	7.33	70	248972	37.189	ng	98
20) 3+4-Methylphenols	7.32	107	384810	38.619	ng	89
22) Acetophenone	7.33	105	520613	38.929	ng	# 96
24) Nitrobenzene	7.50	77	398649	40.830	ng	98
25) Isophorone	7.74	82	682816	38.371	ng	99
26) 2-Nitrophenol	7.82	139	169433	36.087	ng	91
27) 2,4-Dimethylphenol	7.85	122	340869	38.443	ng	98
28) bis(2-Chloroethoxy)methane	7.95	93	433817	39.119	ng	100
29) 2,4-Dichlorophenol	8.05	162	298468	39.206	ng	100
30) 1,2,4-Trichlorobenzene	8.14	180	297344	39.052	ng	99
31) Naphthalene	8.22	128	1011307	39.010	ng	99
32) Benzoic acid	7.95	122	163943	22.509	ng	98
33) 4-Chloroaniline	8.27	127	424501	39.212	ng	97
34) Hexachlorobutadiene	8.33	225	153235	39.073	ng	98
35) Caprolactam	8.65	113	92495	37.877	ng	96
36) 4-Chloro-3-methylphenol	8.74	107	324902	37.971	ng	99
37) 2-Methylnaphthalene	8.91	142	637119	37.805	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	287783	42.124	ng	99
40) Hexachlorocyclopentadiene	9.06	237	15933	5.103	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.19	196	192293	39.731	ng	99
43) 2,4,5-Trichlorophenol	9.23	196	190185	39.365	ng	96
45) 1,1'-Biphenyl	9.37	154	822867	41.796	ng	98
46) 2-Chloronaphthalene	9.40	162	596903	40.380	ng	97
47) 2-Nitroaniline	9.50	65	199156	44.000	ng	92
48) Acenaphthylene	9.82	152	901394	39.834	ng	99
49) Dimethylphthalate	9.67	163	709818	41.055	ng	99
50) 2,6-Dinitrotoluene	9.74	165	151078	40.137	ng	90
51) Acenaphthene	9.99	154	534771	37.307	ng	99
52) 3-Nitroaniline	9.90	138	188865	43.033	ng	98
53) 2,4-Dinitrophenol	10.01	184	4462	7.939	ng	94
54) Dibenzofuran	10.16	168	749960	39.295	ng	99
55) 4-Nitrophenol	10.06	139	128841	34.718	ng	93
56) 2,4-Dinitrotoluene	10.14	165	186322	38.141	ng	94
57) Fluorene	10.50	166	590089	38.467	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.27	232	119476	35.798	ng	97
59) Diethylphthalate	10.37	149	678806	40.179	ng	99
60) 4-Chlorophenyl-phenylether	10.49	204	267708	38.850	ng	99
61) 4-Nitroaniline	10.52	138	179459	42.383	ng	96
62) Azobenzene	10.65	77	593356	38.197	ng	99
64) 4,6-Dinitro-2-methylphenol	10.55	198	11216	5.031	ng	89
65) n-Nitrosodiphenylamine	10.61	169	544585	42.901	ng	99
66) 4-Bromophenyl-phenylether	10.99	248	148451	41.390	ng	92
67) Hexachlorobenzene	11.05	284	152473	39.803	ng	97
68) Atrazine	11.13	200	142642	37.349	ng	99
69) Pentachlorophenol	11.25	266	82886	34.766	ng	99
70) Phenanthrene	11.47	178	770147	39.266	ng	99
71) Anthracene	11.52	178	797204	39.724	ng	100
72) Carbazole	11.67	167	782521	39.818	ng	100
73) Di-n-butylphthalate	12.00	149	963971	40.751	ng	99
74) Fluoranthene	12.66	202	775104	39.027	ng	99
76) Benzidine	12.77	184	473916	45.727	ng	100
77) Pyrene	12.89	202	804046	37.630	ng	100
79) Butylbenzylphthalate	13.50	149	403629	40.194	ng	96
80) Benzo(a)anthracene	14.07	228	666948	39.307	ng	99
81) 3,3'-Dichlorobenzidine	14.03	252	266318	42.816	ng	99
82) Chrysene	14.11	228	638231	39.509	ng	99
83) Bis(2-ethylhexyl)phthalate	14.05	149	570158	41.234	ng	# 99
84) Di-n-octyl phthalate	14.67	149	987467	44.350	ng	98
85) Indeno(1,2,3-cd)pyrene	17.09	276	417218	32.287	ng	96
87) Benzo(b)fluoranthene	15.13	252	561906m	42.696	ng	
88) Benzo(k)fluoranthene	15.17	252	532192	40.942	ng	98
89) Benzo(a)pyrene	15.51	252	477653	39.539	ng	# 96
90) Dibenzo(a,h)anthracene	17.10	278	354575	36.675	ng	98
91) Benzo(g,h,i)perylene	17.54	276	326988	34.216	ng	95

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

