

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097026.D
 Acq On : 25 Jul 2017 13:56
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:54:19 PM

Quant Time: Jul 25 15:01:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	152481	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	597781	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	256637	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	414507	20.00	ng	0.00
75) Chrysene-d12	13.63	240	270178	20.00	ng	0.00
86) Perylene-d12	14.97	264	273342	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.10	112	756755	74.82	ng	0.00
7) Phenol-d6	6.17	99	889583	77.04	ng	0.00
23) Nitrobenzene-d5	7.07	82	849002	83.32	ng	0.00
41) 2,4,6-Tribromophenol	10.32	330	154618	76.25	ng	0.00
44) 2-Fluorobiphenyl	8.87	172	1161759	76.83	ng	0.00
78) Terphenyl-d14	12.59	244	1015550	76.64	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.04	88	149865	35.504	ng	# 74
3) Pyridine	2.66	79	481436	37.248	ng	# 69
4) n-Nitrosodimethylamine	2.62	42	264872	36.990	ng	# 86
6) Aniline	6.16	93	559821	38.526	ng	# 80
8) 2-Chlorophenol	6.29	128	422484	39.062	ng	# 95
9) Benzaldehyde	6.05	77	272226	39.828	ng	# 98
10) Phenol	6.18	94	549923	40.149	ng	# 96
11) bis(2-Chloroethyl)ether	6.25	93	440602	40.672	ng	# 90
12) 1,3-Dichlorobenzene	6.44	146	440710	37.811	ng	# 97
13) 1,4-Dichlorobenzene	6.52	146	443289m	38.377	ng	# 97
14) 1,2-Dichlorobenzene	6.67	146	383085	37.983	ng	# 97
15) Benzyl Alcohol	6.66	79	273550	37.417	ng	# 96
16) 2,2'-oxybis(1-Chloropropan	6.79	45	857571	39.469	ng	# 84
17) 2-Methylphenol	6.79	107	318011	38.097	ng	# 91
18) Hexachloroethane	7.01	117	166480	39.395	ng	# 79
19) n-Nitroso-di-n-propylamine	6.93	70	295731	38.908	ng	# 83
20) 3+4-Methylphenols	6.95	107	423415	38.837	ng	# 62
22) Acetophenone	6.92	105	579052	40.337	ng	# 99
24) Nitrobenzene	7.09	77	445529	41.981	ng	# 96
25) Isophorone	7.34	82	757817	37.975	ng	# 97
26) 2-Nitrophenol	7.41	139	227838	39.682	ng	# 90
27) 2,4-Dimethylphenol	7.47	122	353251	37.297	ng	# 95
28) bis(2-Chloroethoxy)methane	7.55	93	483732	37.145	ng	# 98
29) 2,4-Dichlorophenol	7.66	162	322906	39.575	ng	# 96
30) 1,2,4-Trichlorobenzene	7.73	180	289066	37.168	ng	# 96
31) Naphthalene	7.81	128	1062119	37.692	ng	# 99
32) Benzoic acid	7.63	122	271638	38.408	ng	# 93
33) 4-Chloroaniline	7.86	127	438717	38.263	ng	# 99
34) Hexachlorobutadiene	7.93	225	155616	37.743	ng	# 97
35) Caprolactam	8.25	113	101667	38.858	ng	# 47
36) 4-Chloro-3-methylphenol	8.37	107	348972	39.203	ng	# 89
37) 2-Methylnaphthalene	8.50	142	655592	36.746	ng	# 99
39) 1,2,4,5-Tetrachlorobenzene	8.67	216	261465	38.183	ng	# 100
40) Hexachlorocyclopentadiene	8.66	237	81032	37.643	ng	# 98

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097026.D
 Acq On : 25 Jul 2017 13:56
 Operator : SJ/JU
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDC0040

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:54:19 PM

Quant Time: Jul 25 15:01:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.79	196	192963	38.885	nq	99
43) 2,4,5-Trichlorophenol	8.83	196	191348	38.524	nq	99
45) 1,1'-Biphenyl	8.96	154	822990	37.545	nq	98
46) 2-Chloronaphthalene	8.99	162	605128	37.514	nq	94
47) 2-Nitroaniline	9.09	65	244428	41.753	nq	98
48) Acenaphthylene	9.39	152	938338	37.898	nq	99
49) Dimethylphthalate	9.27	163	743903	38.171	nq	99
50) 2,6-Dinitrotoluene	9.33	165	161532	39.080	nq	# 81
51) Acenaphthene	9.56	154	554220	38.001	nq	100
52) 3-Nitroaniline	9.50	138	200770	39.132	nq	94
53) 2,4-Dinitrophenol	9.61	184	77326	41.145	nq	# 74
54) Dibenzofuran	9.74	168	712174	36.661	nq	99
55) 4-Nitrophenol	9.68	139	118825	38.945	nq	# 63
56) 2,4-Dinitrotoluene	9.73	165	178495	37.565	nq	# 48
57) Fluorene	10.08	166	577911	39.582	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.86	232	129337	37.713	nq	100
59) Diethylphthalate	9.97	149	722973	40.436	nq	100
60) 4-Chlorophenyl-phenylether	10.07	204	238105	39.492	nq	95
61) 4-Nitroaniline	10.11	138	201395	41.312	nq	92
62) Azobenzene	10.23	77	643159	38.776	nq	93
64) 4,6-Dinitro-2-methylphenol	10.15	198	114618	44.500	nq	93
65) n-Nitrosodiphenylamine	10.20	169	558581	38.730	nq	99
66) 4-Bromophenyl-phenylether	10.56	248	163439	39.510	nq	95
67) Hexachlorobenzene	10.63	284	177244	38.209	nq	# 90
68) Atrazine	10.73	200	158438	38.310	nq	92
69) Pentachlorophenol	10.82	266	74340	38.732	nq	98
70) Phenanthrene	11.03	178	840112	37.827	nq	99
71) Anthracene	11.08	178	846333	37.206	nq	99
72) Carbazole	11.24	167	851190	38.048	nq	100
73) Di-n-butylphthalate	11.58	149	1084158	39.205	nq	99
74) Fluoranthene	12.21	202	826420	37.983	nq	99
76) Benzidine	12.33	184	419295	41.825	nq	97
77) Pyrene	12.43	202	841914	38.064	nq	99
79) Butylbenzylphthalate	13.07	149	451040	40.088	nq	# 82
80) Benzo(a)anthracene	13.62	228	677547	38.758	nq	99
81) 3,3'-Dichlorobenzidine	13.59	252	258305	39.182	nq	97
82) Chrysene	13.66	228	673656	39.464	nq	99
83) Bis(2-ethylhexyl)phthalate	13.63	149	573096	39.012	nq	99
84) Di-n-octyl phthalate	14.24	149	1040086	38.581	nq	# 93
85) Indeno(1,2,3-cd)pyrene	16.20	276	565102	38.607	nq	96
87) Benzo(b)fluoranthene	14.62	252	627920	38.215	nq	98
88) Benzo(k)fluoranthene	14.64	252	586511	37.459	nq	98
89) Benzo(a)pyrene	14.92	252	571286m	37.989	nq	
90) Dibenzo(a,h)anthracene	16.22	278	447781	36.310	nq	97
91) Benzo(g,h,i)perylene	16.57	276	477780	36.945	nq	96

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

