

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050116\
 Data File : BF086593.D
 Acq On : 2 May 2016 2:49
 Operator : UM/SJ
 Sample : H2702-10MSD
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP-B1AMSD

Manual Integrations
 APPROVED

sohil
 5/2/2016 7:48:41 PM

Quant Time: May 02 05:25:33 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 02 04:16:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	146314	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	646556	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	305258	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	585413	20.00	ng	0.00
75) Chrysene-d12	13.92	240	373643	20.00	ng	0.00
86) Perylene-d12	15.30	264	320676	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	1140694	134.47	ng	0.02
7) Phenol-d6	6.42	99	1457099	123.08	ng	0.01
23) Nitrobenzene-d5	7.33	82	1036216	86.39	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	381123	118.39	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	1602301	93.92	ng	-0.01
78) Terphenyl-d14	12.87	244	1375367	81.10	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.49	88	169707	38.09	ng	# 70
3) Pyridine	3.15	79	333135m	31.59	ng	
4) n-Nitrosodimethylamine	3.08	42	295967	49.24	ng	# 58
6) Aniline	6.45	93	104885	7.32	ng	93
8) 2-Chlorophenol	6.56	128	456642	43.22	ng	97
10) Phenol	6.43	94	533379	40.38	ng	97
11) bis(2-Chloroethyl)ether	6.51	93	480361	42.07	ng	90
12) 1,3-Dichlorobenzene	6.70	146	435819m	38.35	ng	
13) 1,4-Dichlorobenzene	6.78	146	470414	40.12	ng	96
14) 1,2-Dichlorobenzene	6.93	146	401619	37.35	ng	95
15) Benzyl Alcohol	6.92	79	321079	42.87	ng	90
16) 2,2'-oxybis(1-Chloropropan	7.05	45	906743	39.67	ng	97
17) 2-Methylphenol	7.04	107	373945	43.79	ng	# 91
18) Hexachloroethane	7.28	117	167045	38.12	ng	98
19) n-Nitroso-di-n-propylamine	7.20	70	350450	44.78	ng	# 87
20) 3+4-Methylphenols	7.18	107	433901	44.50	ng	# 73
22) Acetophenone	7.18	105	660297	46.57	ng	# 94
24) Nitrobenzene	7.36	77	557640	45.19	ng	94
25) Isophorone	7.60	82	950965	41.19	ng	98
26) 2-Nitrophenol	7.66	139	239289	42.12	ng	# 71
27) 2,4-Dimethylphenol	7.72	122	426657	42.80	ng	95
28) bis(2-Chloroethoxy)methane	7.81	93	575041	45.72	ng	99
29) 2,4-Dichlorophenol	7.92	162	380176	45.20	ng	99
30) 1,2,4-Trichlorobenzene	8.00	180	400100	41.01	ng	98
31) Naphthalene	8.08	128	1579268	48.00	ng	99
32) Benzoic acid	7.85	122	201313	36.03	ng	85
33) 4-Chloroaniline	8.17	127	27185	2.21	ng	# 73
34) Hexachlorobutadiene	8.19	225	204752	37.43	ng	98
35) Caprolactam	8.50	113	92683m	33.03	ng	
36) 4-Chloro-3-methylphenol	8.62	107	419285	43.54	ng	95
37) 2-Methylnaphthalene	8.76	142	824801	42.49	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.93	216	360029	41.20	ng	97
40) Hexachlorocyclopentadiene	8.92	237	364931	84.05	ng	95
42) 2,4,6-Trichlorophenol	9.05	196	230950	41.81	ng	93

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050116\
 Data File : BF086593.D
 Acq On : 2 May 2016 2:49
 Operator : UM/SJ
 Sample : H2702-10MSD
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP-B1AMSD

Manual Integrations
 APPROVED

sohil
 5/2/2016 7:48:41 PM

Quant Time: May 02 05:25:33 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 02 04:16:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.09	196	241410	40.10	ng	# 86
45) 1,1'-Biphenyl	9.23	154	1053998	41.31	ng	97
46) 2-Chloronaphthalene	9.25	162	799563	41.21	ng	94
47) 2-Nitroaniline	9.36	65	284677	45.76	ng	91
48) Acenaphthylene	9.66	152	1166639	38.47	ng	99
49) Dimethylphthalate	9.54	163	962677	41.89	ng	100
50) 2,6-Dinitrotoluene	9.60	165	207148	43.63	ng	92
51) Acenaphthene	9.84	154	866893	44.51	ng	99
52) 3-Nitroaniline	9.77	138	46745	8.66	ng	93
53) 2,4-Dinitrophenol	9.87	184	189999	73.86	ng	# 78
54) Dibenzofuran	10.01	168	1025783	42.14	ng	94
55) 4-Nitrophenol	9.94	139	220728	62.44	ng	81
56) 2,4-Dinitrotoluene	10.00	165	253264	41.85	ng	# 83
57) Fluorene	10.35	166	782368	37.01	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.13	232	218507	46.43	ng	96
59) Diethylphthalate	10.24	149	907764	41.73	ng	98
60) 4-Chlorophenyl-phenylether	10.35	204	384919	40.11	ng	# 82
61) 4-Nitroaniline	10.37	138	164304	31.13	ng	# 77
62) Azobenzene	10.50	77	1003060	42.09	ng	84
64) 4,6-Dinitro-2-methylphenol	10.41	198	146578	42.65	ng	90
65) n-Nitrosodiphenylamine	10.46	169	602821	32.95	ng	100
66) 4-Bromophenyl-phenylether	10.83	248	238004	39.45	ng	95
67) Hexachlorobenzene	10.90	284	284717	40.78	ng	# 93
68) Atrazine	10.99	200	114240	20.79	ng	93
69) Pentachlorophenol	11.09	266	287315	78.05	ng	97
70) Phenanthrene	11.31	178	1231197	40.08	ng	99
71) Anthracene	11.36	178	1255481	38.29	ng	100
72) Carbazole	11.52	167	1001145	34.45	ng	99
73) Di-n-butylphthalate	11.85	149	1267811	38.51	ng	# 98
74) Fluoranthene	12.49	202	1161084	37.44	ng	99
77) Pyrene	12.72	202	1174458	43.62	ng	100
79) Butylbenzylphthalate	13.35	149	539019	46.23	ng	# 86
80) Benzo(a)anthracene	13.91	228	776575	38.98	ng	99
81) 3,3'-Dichlorobenzidine	13.87	252	38241	2.87	ng	95
82) Chrysene	13.94	228	929259	42.92	ng	100
83) Bis(2-ethylhexyl)phthalate	13.91	149	599280	41.40	ng	98
84) Di-n-octyl phthalate	14.51	149	937499	41.36	ng	# 87
85) Indeno(1,2,3-cd)pyrene	16.64	276	840178	52.37	ng	96
87) Benzo(b)fluoranthene	14.91	252	943087m	49.99	ng	
88) Benzo(k)fluoranthene	14.95	252	627491m	31.89	ng	
89) Benzo(a)pyrene	15.25	252	762379	42.80	ng	98
90) Dibenzo(a,h)anthracene	16.66	278	703289	48.25	ng	# 95
91) Benzo(g,h,i)perylene	17.04	276	689854	47.23	ng	98

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

