

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
 Data File : BF085642.D
 Acq On : 22 Mar 2016 2:31
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/22/2016 5:34:35 PM

Quant Time: Mar 22 05:23:39 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	180832	20.00	ng	-0.01
21) Naphthalene-d8	8.13	136	678249	20.00	ng	-0.01
38) Acenaphthene-d10	9.89	164	296661	20.00	ng	-0.01
63) Phenanthrene-d10	11.36	188	447841	20.00	ng	-0.02
75) Chrysene-d12	14.00	240	344151	20.00	ng	-0.01
86) Perylene-d12	15.42	264	345450	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	891405	82.80	ng	-0.01
7) Phenol-d6	6.48	99	1027744	74.43	ng	-0.01
23) Nitrobenzene-d5	7.42	82	934764	80.00	ng	-0.01
41) 2,4,6-Tribromophenol	10.68	330	217615	74.89	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	1552980	90.87	ng	-0.01
78) Terphenyl-d14	12.95	244	1067995	74.68	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.42	88	211619	40.44	ng	100
3) Pyridine	3.14	79	569431	44.31	ng	98
4) n-Nitrosodimethylamine	3.11	42	207754	38.73	ng	99
6) Aniline	6.50	93	722410	38.95	ng	# 80
8) 2-Chlorophenol	6.63	128	485626	38.99	ng	97
9) Benzaldehyde	6.39	77	265967	41.35	ng	100
10) Phenol	6.49	94	535244	34.03	ng	85
11) bis(2-Chloroethyl)ether	6.58	93	446200	37.63	ng	98
12) 1,3-Dichlorobenzene	6.79	146	547212	40.32	ng	100
13) 1,4-Dichlorobenzene	6.86	146	515169	37.21	ng	99
14) 1,2-Dichlorobenzene	7.02	146	489519	39.83	ng	99
15) Benzyl Alcohol	7.00	79	366485	38.75	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.12	45	615263	39.46	ng	95
17) 2-Methylphenol	7.11	107	418461	40.53	ng	99
18) Hexachloroethane	7.36	117	195235	39.06	ng	94
19) n-Nitroso-di-n-propylamine	7.27	70	296626	35.96	ng	94
20) 3+4-Methylphenols	7.26	107	412154	34.47	ng	94
22) Acetophenone	7.26	105	605407	37.61	ng	# 100
24) Nitrobenzene	7.43	77	491259	38.41	ng	97
25) Isophorone	7.67	82	957638	40.94	ng	100
26) 2-Nitrophenol	7.75	139	278269	44.02	ng	98
27) 2,4-Dimethylphenol	7.80	122	453814	40.60	ng	95
28) bis(2-Chloroethoxy)methane	7.89	93	572712	43.04	ng	99
29) 2,4-Dichlorophenol	7.99	162	347545	37.65	ng	98
30) 1,2,4-Trichlorobenzene	8.07	180	400801	38.76	ng	99
31) Naphthalene	8.15	128	1218199	35.58	ng	100
32) Benzoic acid	7.93	122	277101m	29.65	ng	
33) 4-Chloroaniline	8.21	127	567906	40.47	ng	99
34) Hexachlorobutadiene	8.28	225	223430	40.55	ng	99
35) Caprolactam	8.60	113	125042	39.37	ng	92
36) 4-Chloro-3-methylphenol	8.69	107	378893	35.14	ng	96
37) 2-Methylnaphthalene	8.85	142	844765	40.48	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	348297	38.70	ng	99
40) Hexachlorocyclopentadiene	9.00	237	69236	17.07	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.12	196	234255	38.03	ng	98
43) 2,4,5-Trichlorophenol	9.17	196	255420	40.74	ng	96
45) 1,1'-Biphenyl	9.30	154	959333	37.49	ng	98
46) 2-Chloronaphthalene	9.34	162	770890	41.24	ng	98
47) 2-Nitroaniline	9.43	65	233539	41.30	ng	100
48) Acenaphthylene	9.75	152	1212131	40.07	ng	100
49) Dimethylphthalate	9.61	163	879041	39.27	ng	100
50) 2,6-Dinitrotoluene	9.67	165	176804	37.41	ng	95
51) Acenaphthene	9.92	154	728850	38.39	ng	99
52) 3-Nitroaniline	9.84	138	204244	34.51	ng	100
53) 2,4-Dinitrophenol	9.94	184	35312	12.20	ng	91
54) Dibenzofuran	10.09	168	918287	39.72	ng	96
55) 4-Nitrophenol	10.00	139	135995	29.73	ng	94
56) 2,4-Dinitrotoluene	10.08	165	244331	39.49	ng	94
57) Fluorene	10.44	166	737068	38.25	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.21	232	158842	35.19	ng	98
59) Diethylphthalate	10.31	149	899352	40.41	ng	99
60) 4-Chlorophenyl-phenylether	10.42	204	363255	38.61	ng	92
61) 4-Nitroaniline	10.46	138	215468	36.93	ng	96
62) Azobenzene	10.58	77	814000	38.11	ng	96
64) 4,6-Dinitro-2-methylphenol	10.48	198	63111	22.02	ng	99
65) n-Nitrosodiphenylamine	10.54	169	637625	45.71	ng	99
66) 4-Bromophenyl-phenylether	10.92	248	225196	50.34	ng	# 91
67) Hexachlorobenzene	10.98	284	215957	45.44	ng	# 89
68) Atrazine	11.08	200	182600	42.99	ng	94
69) Pentachlorophenol	11.17	266	100060	34.67	ng	99
70) Phenanthrene	11.40	178	966875	40.95	ng	99
71) Anthracene	11.44	178	923748	37.31	ng	100
72) Carbazole	11.60	167	821753	38.47	ng	97
73) Di-n-butylphthalate	11.93	149	1158530	43.31	ng	98
74) Fluoranthene	12.57	202	795294	32.17	ng	97
76) Benzidine	12.70	184	378579	32.71	ng	98
77) Pyrene	12.80	202	798071	32.30	ng	99
79) Butylbenzylphthalate	13.42	149	385801	32.75	ng	# 76
80) Benzo(a)anthracene	13.99	228	759959	38.04	ng	99
81) 3,3'-Dichlorobenzidine	13.96	252	299555	40.91	ng	# 96
82) Chrysene	14.03	228	639545	33.27	ng	100
83) Bis(2-ethylhexyl)phthalate	13.98	149	501675	34.28	ng	# 98
84) Di-n-octyl phthalate	14.60	149	985019	41.31	ng	99
85) Indeno(1,2,3-cd)pyrene	16.83	276	789124	49.13	ng	99
87) Benzo(b)fluoranthene	15.02	252	829209	36.79	ng	# 97
88) Benzo(k)fluoranthene	15.04	252	735513m	39.61	ng	
89) Benzo(a)pyrene	15.36	252	776335	40.63	ng	98
90) Dibenzo(a,h)anthracene	16.84	278	676095	42.42	ng	99
91) Benzo(g,h,i)perylene	17.24	276	607315	39.37	ng	98

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

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