

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072215\
 Data File : BF080587.D
 Acq On : 23 Jul 2015 7:22
 Operator : TP/UM
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 7/23/2015 11:51:53 AM

Quant Time: Jul 23 07:48:27 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 23 05:13:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	34967	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	137078	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	61415	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	105417	20.00	ng	0.00
75) Chrysene-d12	14.34	240	68011	20.00	ng	0.02
86) Perylene-d12	16.07	264	39613	20.00	ng	0.05

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.53	112	161703	82.01	ng	0.00
7) Phenol-d6	6.57	99	208122	85.70	ng	0.00
23) Nitrobenzene-d5	7.52	82	206322	96.95	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	40026	89.57	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	347445	81.35	ng	0.00
78) Terphenyl-d14	13.14	244	315639	95.17	ng	0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.50	88	35758	34.89	ng	# 100
3) Pyridine	3.28	79	99945	42.00	ng	94
4) n-Nitrosodimethylamine	3.18	42	64519	42.92	ng	# 93
6) Aniline	6.61	93	146899	40.37	ng	95
8) 2-Chlorophenol	6.73	128	88559	42.89	ng	# 71
9) Benzaldehyde	6.50	77	72432	39.64	ng	93
10) Phenol	6.58	94	129218	44.74	ng	99
11) bis(2-Chloroethyl)ether	6.68	93	89415	43.10	ng	94
12) 1,3-Dichlorobenzene	6.90	146	109998	41.18	ng	99
13) 1,4-Dichlorobenzene	6.98	146	104400	41.03	ng	100
14) 1,2-Dichlorobenzene	7.13	146	106610	43.44	ng	99
15) Benzyl Alcohol	7.09	79	86541	43.04	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.24	45	161875	37.26	ng	97
17) 2-Methylphenol	7.20	107	71561	40.63	ng	94
18) Hexachloroethane	7.48	117	38103	42.48	ng	95
19) n-Nitroso-di-n-propylamine	7.37	70	67048	42.99	ng	# 95
20) 3+4-Methylphenols	7.36	107	93560	42.12	ng	97
22) Acetophenone	7.37	105	130939	40.01	ng	98
24) Nitrobenzene	7.54	77	100848	39.55	ng	99
25) Isophorone	7.78	82	179457	39.92	ng	99
26) 2-Nitrophenol	7.86	139	37724	56.11	ng	95
27) 2,4-Dimethylphenol	7.89	122	79917	42.00	ng	95
28) bis(2-Chloroethoxy)methane	8.00	93	93605	38.82	ng	96
29) 2,4-Dichlorophenol	8.10	162	69580	45.75	ng	95
30) 1,2,4-Trichlorobenzene	8.19	180	85286	40.10	ng	98
31) Naphthalene	8.27	128	278178	40.87	ng	98
32) Benzoic acid	7.95	122	35955m	56.61	ng	
33) 4-Chloroaniline	8.32	127	107744	41.13	ng	97
34) Hexachlorobutadiene	8.40	225	49685	40.00	ng	99
35) Caprolactam	8.65	113	18595	43.35	ng	# 65
36) 4-Chloro-3-methylphenol	8.78	107	83711	44.51	ng	99
37) 2-Methylnaphthalene	8.96	142	157434	41.30	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	87061	41.35	ng	99
40) Hexachlorocyclopentadiene	9.12	237	45360	50.00	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	47474	45.53	ng	100
43) 2,4,5-Trichlorophenol	9.26	196	50096	45.97	ng	98
45) 1,1'-Biphenyl	9.42	154	205286	39.88	ng	99
46) 2-Chloronaphthalene	9.45	162	167888	41.29	ng	96
47) 2-Nitroaniline	9.54	65	46686	48.67	ng	92
48) Acenaphthylene	9.87	152	243152	40.74	ng	99
49) Dimethylphthalate	9.72	163	188776	43.77	ng	99
50) 2,6-Dinitrotoluene	9.78	165	34881	48.24	ng	97
51) Acenaphthene	10.04	154	142364	40.61	ng	100
52) 3-Nitroaniline	9.95	138	33889	46.23	ng	97
53) 2,4-Dinitrophenol	10.05	184	7789	60.55	ng #	73
54) Dibenzofuran	10.21	168	209137	40.61	ng	99
55) 4-Nitrophenol	10.09	139	26916	47.46	ng	100
56) 2,4-Dinitrotoluene	10.18	165	40889	49.63	ng	96
57) Fluorene	10.56	166	170171	41.35	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.33	232	37230	42.20	ng	97
59) Diethylphthalate	10.42	149	170661	42.88	ng	98
60) 4-Chlorophenyl-phenylether	10.54	204	74483	41.97	ng	99
61) 4-Nitroaniline	10.56	138	34480	43.28	ng	95
62) Azobenzene	10.70	77	184676	42.13	ng	100
64) 4,6-Dinitro-2-methylphenol	10.59	198	11759	56.14	ng	88
65) n-Nitrosodiphenylamine	10.66	169	145408	41.84	ng	98
66) 4-Bromophenyl-phenylether	11.04	248	45277	44.63	ng	94
67) Hexachlorobenzene	11.10	284	49341	42.81	ng	100
68) Atrazine	11.18	200	44104	47.30	ng	98
69) Pentachlorophenol	11.29	266	22810	49.85	ng	99
70) Phenanthrene	11.52	178	240718	40.23	ng	98
71) Anthracene	11.57	178	220245	39.77	ng	99
72) Carbazole	11.72	167	200205	39.91	ng	99
73) Di-n-butylphthalate	12.06	149	260995	44.61	ng #	98
74) Fluoranthene	12.74	202	216485	39.99	ng	97
76) Benzidine	12.87	184	98523	48.72	ng	97
77) Pyrene	12.98	202	214894	46.41	ng	97
79) Butylbenzylphthalate	13.68	149	74435	49.71	ng #	78
80) Benzo(a)anthracene	14.33	228	160610	41.56	ng	99
81) 3,3'-Dichlorobenzidine	14.29	252	47016	40.27	ng	95
82) Chrysene	14.37	228	153942	39.86	ng	96
83) Bis(2-ethylhexyl)phthalate	14.35	149	107283	43.25	ng	97
84) Di-n-octyl phthalate	15.13	149	135873	43.78	ng	96
85) Indeno(1,2,3-cd)pyrene	17.60	276	157880	57.53	ng	98
87) Benzo(b)fluoranthene	15.61	252	96795	38.86	ng	98
88) Benzo(k)fluoranthene	15.64	252	111027m	45.02	ng	
89) Benzo(a)pyrene	16.00	252	89716	41.57	ng #	93
90) Dibenzo(a,h)anthracene	17.62	278	123355	66.96	ng	99
91) Benzo(g,h,i)perylene	18.05	276	168734	90.07	ng #	92

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

