

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022615\
 Data File : BF077501.D
 Acq On : 27 Feb 2015 10:13
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 2/27/2015 4:21:48 PM

Quant Time: Feb 27 12:12:11 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 27 07:02:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.27	152	57792	20.00	ng	0.00
21) Naphthalene-d8	8.85	136	243363	20.00	ng	0.00
38) Acenaphthene-d10	11.02	164	122104	20.00	ng	0.00
63) Phenanthrene-d10	12.86	188	243938	20.00	ng	0.00
75) Chrysene-d12	16.16	240	268955	20.00	ng	0.02
86) Perylene-d12	17.91	264	265638	20.00	ng	0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	324089	93.62	ng	0.00
7) Phenol-d6	6.83	99	378612	85.06	ng	0.00
23) Nitrobenzene-d5	7.96	82	311333	79.55	ng	0.00
41) 2,4,6-Tribromophenol	12.00	330	100810	85.74	ng	0.00
44) 2-Fluorobiphenyl	10.19	172	651898	83.25	ng	0.00
78) Terphenyl-d14	14.86	244	856364	74.44	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.20	88	64000	43.57	ng	94
3) Pyridine	2.95	79	188962	44.96	ng	96
4) n-Nitrosodimethylamine	2.86	42	73457	40.20	ng	95
6) Aniline	6.85	93	246068	40.00	ng	# 90
8) 2-Chlorophenol	7.00	128	159274	39.00	ng	90
9) Benzaldehyde	6.72	77	137127	50.75	ng	95
10) Phenol	6.84	94	214908	40.69	ng	91
11) bis(2-Chloroethyl)ether	6.96	93	141847	37.38	ng	90
12) 1,3-Dichlorobenzene	7.19	146	173888	39.10	ng	# 91
13) 1,4-Dichlorobenzene	7.29	146	170346	37.66	ng	96
14) 1,2-Dichlorobenzene	7.47	146	166757	38.75	ng	96
15) Benzyl Alcohol	7.45	79	129561	38.29	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.63	45	206089	37.99	ng	98
17) 2-Methylphenol	7.59	107	127605	39.98	ng	99
18) Hexachloroethane	7.89	117	53000	35.08	ng	91
19) n-Nitroso-di-n-propylamine	7.78	70	107066	37.88	ng	94
20) 3+4-Methylphenols	7.79	107	165762	39.43	ng	# 76
22) Acetophenone	7.78	105	220879	38.58	ng	# 89
24) Nitrobenzene	7.98	77	163883	39.80	ng	89
25) Isophorone	8.28	82	288426	37.15	ng	96
26) 2-Nitrophenol	8.38	139	63553	37.66	ng	# 87
27) 2,4-Dimethylphenol	8.44	122	142856	37.84	ng	98
28) bis(2-Chloroethoxy)methane	8.56	93	178483	36.39	ng	98
29) 2,4-Dichlorophenol	8.67	162	134922	38.07	ng	88
30) 1,2,4-Trichlorobenzene	8.78	180	143489	36.82	ng	96
31) Naphthalene	8.87	128	474891	39.02	ng	99
32) Benzoic acid	8.53	122	73685	40.16	ng	99
33) 4-Chloroaniline	8.94	127	205909	38.05	ng	# 92
34) Hexachlorobutadiene	9.03	225	82560	37.11	ng	99
35) Caprolactam	9.37	113	41933	38.76	ng	95
36) 4-Chloro-3-methylphenol	9.55	107	141047	39.59	ng	100
37) 2-Methylnaphthalene	9.73	142	331414	38.35	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.94	216	149353	42.63	ng	99
40) Hexachlorocyclopentadiene	9.92	237	19198	10.22	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.08	196	99341	42.82	ng	98
43) 2,4,5-Trichlorophenol	10.13	196	103386	41.67	ng	96
45) 1,1'-Biphenyl	10.31	154	399503	43.19	ng	99
46) 2-Chloronaphthalene	10.34	162	307673	42.41	ng	94
47) 2-Nitroaniline	10.46	65	85661	47.97	ng	# 70
48) Acenaphthylene	10.85	152	487648	42.46	ng	100
49) Dimethylphthalate	10.70	163	352822	41.11	ng	98
50) 2,6-Dinitrotoluene	10.77	165	75131	43.65	ng	# 74
51) Acenaphthene	11.06	154	275422	40.19	ng	99
52) 3-Nitroaniline	10.97	138	91801	46.26	ng	# 77
53) 2,4-Dinitrophenol	11.11	184	2886	5.51	ng	# 71
54) Dibenzofuran	11.28	168	431179	40.75	ng	98
55) 4-Nitrophenol	11.19	139	66816	41.86	ng	# 72
56) 2,4-Dinitrotoluene	11.27	165	89840	38.63	ng	# 71
57) Fluorene	11.70	166	352061	41.61	ng	98
58) 2,3,4,6-Tetrachlorophenol	11.43	232	84789	42.51	ng	97
59) Diethylphthalate	11.58	149	351299	41.52	ng	95
60) 4-Chlorophenyl-phenylether	11.71	204	160408	39.35	ng	90
61) 4-Nitroaniline	11.73	138	91078	47.88	ng	95
62) Azobenzene	11.91	77	325985	40.61	ng	97
64) 4,6-Dinitro-2-methylphenol	11.77	198	6435	9.81	ng	# 47
65) n-Nitrosodiphenylamine	11.86	169	313598	38.75	ng	99
66) 4-Bromophenyl-phenylether	12.32	248	101590	35.56	ng	# 90
67) Hexachlorobenzene	12.38	284	109833	35.82	ng	# 77
68) Atrazine	12.53	200	95802	36.41	ng	96
69) Pentachlorophenol	12.63	266	56290	34.93	ng	97
70) Phenanthrene	12.90	178	542240	38.35	ng	100
71) Anthracene	12.96	178	525448	37.45	ng	99
72) Carbazole	13.16	167	500914	37.55	ng	99
73) Di-n-butylphthalate	13.62	149	600475	38.33	ng	98
74) Fluoranthene	14.37	202	602890	39.04	ng	100
76) Benzidine	14.55	184	445472	49.02	ng	97
77) Pyrene	14.65	202	614532	37.35	ng	99
79) Butylbenzylphthalate	15.48	149	275386	40.69	ng	98
80) Benzo(a)anthracene	16.15	228	587422	37.27	ng	99
81) 3,3'-Dichlorobenzidine	16.12	252	235674	40.80	ng	# 96
82) Chrysene	16.19	228	561427	37.93	ng	99
83) Bis(2-ethylhexyl)phthalate	16.20	149	410995	37.87	ng	# 95
84) Di-n-octyl phthalate	17.01	149	714416	39.42	ng	99
85) Indeno(1,2,3-cd)pyrene	19.38	276	610139m	36.15	ng	
87) Benzo(b)fluoranthene	17.46	252	676215	43.78	ng	# 96
88) Benzo(k)fluoranthene	17.49	252	482069m	31.84	ng	
89) Benzo(a)pyrene	17.85	252	557644	37.90	ng	99
90) Dibenzo(a,h)anthracene	19.41	278	503457	35.88	ng	99
91) Benzo(g,h,i)perylene	19.81	276	499793	35.29	ng	98

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

