

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
 Data File : BF085828.D
 Acq On : 29 Mar 2016 3:49
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/29/2016 12:03:42 PM

Quant Time: Mar 29 04:55:41 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 15:27:05 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	124557	20.00	ng	-0.01
21) Naphthalene-d8	8.10	136	507087	20.00	ng	-0.01
38) Acenaphthene-d10	9.86	164	216329	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	377093	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	246163	20.00	ng	0.00
86) Perylene-d12	15.40	264	246417	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	600300	80.26	ng	0.00
7) Phenol-d6	6.46	99	777190	81.73	ng	0.00
23) Nitrobenzene-d5	7.38	82	687576	76.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	157331	76.55	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	1074883	83.02	ng	-0.01
78) Terphenyl-d14	12.93	244	957628	80.01	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.37	88	152003	42.56	ng	99
3) Pyridine	3.09	79	395485	46.19	ng	98
4) n-Nitrosodimethylamine	3.04	42	153043	44.65	ng	99
6) Aniline	6.48	93	521871	40.77	ng	99
8) 2-Chlorophenol	6.61	128	350425	40.33	ng	92
9) Benzaldehyde	6.37	77	219219	38.26	ng	91
10) Phenol	6.47	94	467835	43.43	ng	91
11) bis(2-Chloroethyl)ether	6.56	93	343195	41.40	ng	94
12) 1,3-Dichlorobenzene	6.76	146	383438	40.98	ng	# 93
13) 1,4-Dichlorobenzene	6.84	146	399437	42.09	ng	95
14) 1,2-Dichlorobenzene	6.98	146	335665	37.95	ng	94
15) Benzyl Alcohol	6.96	79	242398	38.21	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.10	45	421339	38.30	ng	96
17) 2-Methylphenol	7.08	107	281208	40.42	ng	99
18) Hexachloroethane	7.33	117	140209	40.35	ng	# 86
19) n-Nitroso-di-n-propylamine	7.24	70	202025	35.48	ng	92
20) 3+4-Methylphenols	7.24	107	308486	36.88	ng	97
22) Acetophenone	7.24	105	462868	39.18	ng	# 96
24) Nitrobenzene	7.41	77	400500	43.15	ng	96
25) Isophorone	7.65	82	673030	38.80	ng	97
26) 2-Nitrophenol	7.73	139	193839	41.74	ng	91
27) 2,4-Dimethylphenol	7.76	122	300472	37.02	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	375635	38.00	ng	98
29) 2,4-Dichlorophenol	7.97	162	278831	40.86	ng	99
30) 1,2,4-Trichlorobenzene	8.05	180	300941	39.73	ng	97
31) Naphthalene	8.13	128	1001539	39.20	ng	100
32) Benzoic acid	7.89	122	197305m	36.60	ng	
33) 4-Chloroaniline	8.18	127	400123	38.35	ng	96
34) Hexachlorobutadiene	8.24	225	155878	37.86	ng	99
35) Caprolactam	8.56	113	83895	34.71	ng	100
36) 4-Chloro-3-methylphenol	8.66	107	288114	36.15	ng	98
37) 2-Methylnaphthalene	8.81	142	590456	38.21	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	273687	42.66	ng	99
40) Hexachlorocyclopentadiene	8.97	237	45041	24.10	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	179620	41.46	ng	100
43) 2,4,5-Trichlorophenol	9.14	196	181793	40.01	ng #	94
45) 1,1'-Biphenyl	9.28	154	722114	39.56	ng	99
46) 2-Chloronaphthalene	9.30	162	541342	40.33	ng #	91
47) 2-Nitroaniline	9.41	65	179264	43.67	ng	99
48) Acenaphthylene	9.72	152	884271	40.43	ng	99
49) Dimethylphthalate	9.59	163	711769	43.37	ng	99
50) 2,6-Dinitrotoluene	9.65	165	130942	37.07	ng	94
51) Acenaphthene	9.89	154	531454	39.78	ng	99
52) 3-Nitroaniline	9.82	138	167178	41.34	ng	98
53) 2,4-Dinitrophenol	9.93	184	16856	14.77	ng #	81
54) Dibenzofuran	10.06	168	681009	39.39	ng	94
55) 4-Nitrophenol	9.99	139	87494	33.21	ng	88
56) 2,4-Dinitrotoluene	10.06	165	191745	42.70	ng #	85
57) Fluorene	10.40	166	548380	38.15	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.18	232	115580	36.60	ng	98
59) Diethylphthalate	10.29	149	680309	41.95	ng	98
60) 4-Chlorophenyl-phenylether	10.40	204	303597	43.20	ng #	86
61) 4-Nitroaniline	10.44	138	166245	42.98	ng	95
62) Azobenzene	10.56	77	665203	42.69	ng	91
64) 4,6-Dinitro-2-methylphenol	10.47	198	34718	15.72	ng #	55
65) n-Nitrosodiphenylamine	10.52	169	481979	40.35	ng	98
66) 4-Bromophenyl-phenylether	10.89	248	165979	43.02	ng #	84
67) Hexachlorobenzene	10.96	284	165612	41.05	ng #	86
68) Atrazine	11.05	200	161883	40.15	ng	94
69) Pentachlorophenol	11.16	266	64693	32.66	ng	100
70) Phenanthrene	11.36	178	769021	39.72	ng	100
71) Anthracene	11.42	178	859729	42.30	ng	99
72) Carbazole	11.58	167	728920	42.71	ng	99
73) Di-n-butylphthalate	11.90	149	969292	41.39	ng	99
74) Fluoranthene	12.55	202	692696	33.66	ng	96
76) Benzidine	12.68	184	329350	37.99	ng	99
77) Pyrene	12.78	202	690971	34.15	ng	99
79) Butylbenzylphthalate	13.40	149	333942	39.41	ng #	76
80) Benzo(a)anthracene	13.97	228	586307	41.77	ng	99
81) 3,3'-Dichlorobenzidine	13.93	252	422015	43.32	ng	99
82) Chrysene	14.00	228	476940	33.73	ng	99
83) Bis(2-ethylhexyl)phthalate	13.96	149	444819	43.84	ng #	97
84) Di-n-octyl phthalate	14.56	149	710863	47.12	ng	100
85) Indeno(1,2,3-cd)pyrene	16.79	276	656230	60.11	ng	99
87) Benzo(b)fluoranthene	15.00	252	655297m	42.21	ng	
88) Benzo(k)fluoranthene	15.02	252	449971m	32.15	ng	
89) Benzo(a)pyrene	15.34	252	563562	40.89	ng	98
90) Dibenzo(a,h)anthracene	16.79	278	550701	43.00	ng	98
91) Benzo(g,h,i)perylene	17.20	276	533567	41.15	ng	95

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

