

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
 Data File : BF085802.D
 Acq On : 28 Mar 2016 15:31
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/29/2016 12:03:11 PM

Quant Time: Mar 28 17:08:41 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 15:50:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	123540	20.00	ng	-0.01
21) Naphthalene-d8	8.10	136	518363	20.00	ng	-0.01
38) Acenaphthene-d10	9.86	164	221557	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	401467	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	295641	20.00	ng	0.00
86) Perylene-d12	15.40	264	202605	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	585575	78.94	ng	0.00
7) Phenol-d6	6.46	99	772892	81.95	ng	0.00
23) Nitrobenzene-d5	7.38	82	688907	74.84	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	184558	87.67	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	1118737	84.36	ng	-0.01
78) Terphenyl-d14	12.93	244	1149621	79.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.37	88	144177	40.70	ng	99
3) Pyridine	3.09	79	394184	46.42	ng	98
4) n-Nitrosodimethylamine	3.04	42	143815	42.30	ng	98
6) Aniline	6.48	93	518157	40.81	ng	99
8) 2-Chlorophenol	6.61	128	351457	40.78	ng	91
9) Benzaldehyde	6.37	77	210715	37.08	ng	90
10) Phenol	6.47	94	461863	43.22	ng	88
11) bis(2-Chloroethyl)ether	6.56	93	331020	40.26	ng	93
12) 1,3-Dichlorobenzene	6.76	146	378310	40.76	ng	# 94
13) 1,4-Dichlorobenzene	6.84	146	390302	41.47	ng	95
14) 1,2-Dichlorobenzene	6.98	146	336618	38.37	ng	94
15) Benzyl Alcohol	6.96	79	238567	37.91	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.10	45	419102	38.41	ng	98
17) 2-Methylphenol	7.08	107	278640	40.38	ng	98
18) Hexachloroethane	7.33	117	140499	40.77	ng	# 87
19) n-Nitroso-di-n-propylamine	7.24	70	200109	35.44	ng	93
20) 3+4-Methylphenols	7.24	107	310354	37.41	ng	96
22) Acetophenone	7.24	105	463815	38.41	ng	# 96
24) Nitrobenzene	7.41	77	396416	41.78	ng	96
25) Isophorone	7.65	82	667779	37.66	ng	97
26) 2-Nitrophenol	7.73	139	197394	41.58	ng	91
27) 2,4-Dimethylphenol	7.76	122	297554	35.86	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	372477	36.87	ng	98
29) 2,4-Dichlorophenol	7.97	162	276486	39.64	ng	99
30) 1,2,4-Trichlorobenzene	8.05	180	299555	38.68	ng	98
31) Naphthalene	8.13	128	996663	38.16	ng	100
32) Benzoic acid	7.90	122	233546	42.38	ng	99
33) 4-Chloroaniline	8.18	127	401085	37.61	ng	97
34) Hexachlorobutadiene	8.24	225	152397	36.21	ng	99
35) Caprolactam	8.56	113	83223	33.68	ng	99
36) 4-Chloro-3-methylphenol	8.66	107	294708	36.17	ng	96
37) 2-Methylnaphthalene	8.81	142	596804	37.67	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	271840	41.37	ng	99
40) Hexachlorocyclopentadiene	8.97	237	91774	38.85	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	181972	41.01	ng	100
43) 2,4,5-Trichlorophenol	9.14	196	195184	41.94	ng #	94
45) 1,1'-Biphenyl	9.28	154	717801	38.39	ng	98
46) 2-Chloronaphthalene	9.30	162	544752	39.63	ng #	90
47) 2-Nitroaniline	9.41	65	181718	43.23	ng	96
48) Acenaphthylene	9.72	152	916611	40.92	ng	99
49) Dimethylphthalate	9.59	163	712546	42.39	ng	99
50) 2,6-Dinitrotoluene	9.65	165	134626	37.21	ng	93
51) Acenaphthene	9.90	154	555243	40.58	ng	99
52) 3-Nitroaniline	9.82	138	164192	39.64	ng	98
53) 2,4-Dinitrophenol	9.93	184	71528	42.21	ng #	71
54) Dibenzofuran	10.07	168	709560	40.07	ng	94
55) 4-Nitrophenol	9.99	139	110704	41.03	ng	87
56) 2,4-Dinitrotoluene	10.06	165	202961	44.13	ng #	80
57) Fluorene	10.41	166	589968	40.08	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.18	232	132224	40.88	ng	100
59) Diethylphthalate	10.29	149	709702	42.73	ng	99
60) 4-Chlorophenyl-phenylether	10.40	204	313363	43.54	ng #	88
61) 4-Nitroaniline	10.44	138	180290	45.51	ng	97
62) Azobenzene	10.56	77	704612	44.15	ng	94
64) 4,6-Dinitro-2-methylphenol	10.47	198	106240	45.19	ng #	52
65) n-Nitrosodiphenylamine	10.52	169	506258	39.81	ng	98
66) 4-Bromophenyl-phenylether	10.89	248	178289	43.41	ng #	85
67) Hexachlorobenzene	10.96	284	181651	42.29	ng #	88
68) Atrazine	11.05	200	177670	41.39	ng	96
69) Pentachlorophenol	11.16	266	92579	43.90	ng	99
70) Phenanthrene	11.36	178	887634	43.06	ng	100
71) Anthracene	11.42	178	906352	41.88	ng	99
72) Carbazole	11.58	167	823973	45.35	ng	99
73) Di-n-butylphthalate	11.91	149	1022622	41.02	ng #	98
74) Fluoranthene	12.55	202	822149	37.52	ng	98
76) Benzidine	12.68	184	394973	37.94	ng	99
77) Pyrene	12.78	202	827367	34.05	ng	99
79) Butylbenzylphthalate	13.40	149	373176	36.67	ng #	75
80) Benzo(a)anthracene	13.97	228	665253	39.46	ng	99
81) 3,3'-Dichlorobenzidine	13.93	252	429437	36.71	ng	99
82) Chrysene	14.00	228	563438	33.18	ng	99
83) Bis(2-ethylhexyl)phthalate	13.96	149	464860	38.15	ng #	98
84) Di-n-octyl phthalate	14.56	149	663857	36.64	ng	100
85) Indeno(1,2,3-cd)pyrene	16.78	276	543779	41.47	ng	99
87) Benzo(b)fluoranthene	15.00	252	575767m	45.11	ng	
88) Benzo(k)fluoranthene	15.02	252	398771m	34.66	ng	
89) Benzo(a)pyrene	15.34	252	461584	40.73	ng #	97
90) Dibenzo(a,h)anthracene	16.79	278	454635	43.18	ng	99
91) Benzo(g,h,i)perylene	17.20	276	466829	43.78	ng #	94

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

