

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
 Data File : BF085734.D
 Acq On : 25 Mar 2016 1:23
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/25/2016 3:04:37 PM

Quant Time: Mar 25 02:30:54 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 14:15:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	107309	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	433647	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	212581	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	402688	20.00	ng	0.00
75) Chrysene-d12	13.98	240	286403	20.00	ng	0.00
86) Perylene-d12	15.40	264	199585	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	541969	84.11	ng	0.01
7) Phenol-d6	6.46	99	642288	78.40	ng	0.00
23) Nitrobenzene-d5	7.40	82	621825	80.75	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	163955	81.18	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	1025645	80.61	ng	0.00
78) Terphenyl-d14	12.93	244	991477	71.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.38	88	118598	38.55	ng	99
3) Pyridine	3.11	79	280309	38.00	ng	99
4) n-Nitrosodimethylamine	3.05	42	121774	41.24	ng	99
6) Aniline	6.48	93	436667	39.60	ng	# 69
8) 2-Chlorophenol	6.61	128	309765	41.38	ng	96
9) Benzaldehyde	6.37	77	196766	39.86	ng	94
10) Phenol	6.48	94	345658	37.24	ng	# 56
11) bis(2-Chloroethyl)ether	6.56	93	273841	38.34	ng	96
12) 1,3-Dichlorobenzene	6.77	146	328184	40.71	ng	98
13) 1,4-Dichlorobenzene	6.85	146	328140m	40.13	ng	
14) 1,2-Dichlorobenzene	7.00	146	315162	41.36	ng	98
15) Benzyl Alcohol	6.97	79	225866	41.32	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.11	45	389027	41.05	ng	97
17) 2-Methylphenol	7.09	107	242102	40.39	ng	100
18) Hexachloroethane	7.34	117	122925	41.06	ng	96
19) n-Nitroso-di-n-propylamine	7.25	70	204413	41.67	ng	97
20) 3+4-Methylphenols	7.25	107	295975	41.07	ng	# 86
22) Acetophenone	7.24	105	402915	39.88	ng	99
24) Nitrobenzene	7.42	77	313805	39.53	ng	# 78
25) Isophorone	7.66	82	605079	40.79	ng	# 93
26) 2-Nitrophenol	7.73	139	157511	39.66	ng	# 86
27) 2,4-Dimethylphenol	7.77	122	305540	44.02	ng	96
28) bis(2-Chloroethoxy)methane	7.86	93	347326	41.09	ng	99
29) 2,4-Dichlorophenol	7.98	162	232838	39.90	ng	92
30) 1,2,4-Trichlorobenzene	8.06	180	264017	40.75	ng	93
31) Naphthalene	8.14	128	867625	39.71	ng	98
32) Benzoic acid	7.90	122	212466m	46.09	ng	
33) 4-Chloroaniline	8.18	127	357163	40.03	ng	# 95
34) Hexachlorobutadiene	8.25	225	147008	41.75	ng	98
35) Caprolactam	8.57	113	90652	43.86	ng	# 78
36) 4-Chloro-3-methylphenol	8.68	107	279912	41.07	ng	87
37) 2-Methylnaphthalene	8.82	142	552760	42.90	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	255479	40.53	ng	99
40) Hexachlorocyclopentadiene	8.97	237	82010	36.82	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.11	196	170426	40.03	ng	96
43) 2,4,5-Trichlorophenol	9.14	196	173449	38.85	ng	# 86
45) 1,1'-Biphenyl	9.29	154	745071	41.54	ng	94
46) 2-Chloronaphthalene	9.32	162	561630	42.58	ng	97
47) 2-Nitroaniline	9.41	65	151366	37.53	ng	85
48) Acenaphthylene	9.73	152	871681	40.56	ng	100
49) Dimethylphthalate	9.59	163	611938	37.94	ng	99
50) 2,6-Dinitrotoluene	9.66	165	152715	43.99	ng	# 70
51) Acenaphthene	9.90	154	493140	37.56	ng	99
52) 3-Nitroaniline	9.82	138	161262	40.58	ng	90
53) 2,4-Dinitrophenol	9.93	184	58721	36.99	ng	# 81
54) Dibenzofuran	10.07	168	676649	39.82	ng	98
55) 4-Nitrophenol	9.99	139	103867	40.12	ng	93
56) 2,4-Dinitrotoluene	10.06	165	182983	41.47	ng	# 67
57) Fluorene	10.41	166	579138	41.00	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.20	232	126539	40.78	ng	97
59) Diethylphthalate	10.29	149	612920	38.47	ng	98
60) 4-Chlorophenyl-phenylether	10.40	204	250358	36.25	ng	95
61) 4-Nitroaniline	10.44	138	138812	36.52	ng	87
62) Azobenzene	10.56	77	618845	40.41	ng	97
64) 4,6-Dinitro-2-methylphenol	10.47	198	96750	41.03	ng	# 65
65) n-Nitrosodiphenylamine	10.53	169	525578	41.21	ng	99
66) 4-Bromophenyl-phenylether	10.89	248	168111	40.81	ng	95
67) Hexachlorobenzene	10.96	284	184917	42.92	ng	95
68) Atrazine	11.05	200	165319	38.39	ng	98
69) Pentachlorophenol	11.16	266	88580	41.88	ng	98
70) Phenanthrene	11.37	178	901727	43.61	ng	99
71) Anthracene	11.42	178	859895	39.61	ng	99
72) Carbazole	11.58	167	734760	40.32	ng	99
73) Di-n-butylphthalate	11.91	149	1006760	40.26	ng	100
74) Fluoranthene	12.56	202	862597	39.25	ng	88
76) Benzidine	12.68	184	416556	41.30	ng	99
77) Pyrene	12.78	202	862562	36.64	ng	100
79) Butylbenzylphthalate	13.41	149	390520	39.61	ng	96
80) Benzo(a)anthracene	13.97	228	631642	38.68	ng	99
81) 3,3'-Dichlorobenzidine	13.93	252	407703	35.97	ng	98
82) Chrysene	14.01	228	628350	38.20	ng	99
83) Bis(2-ethylhexyl)phthalate	13.96	149	472099	39.99	ng	99
84) Di-n-octyl phthalate	14.57	149	714253	40.69	ng	100
85) Indeno(1,2,3-cd)pyrene	16.78	276	484896	38.17	ng	99
87) Benzo(b)fluoranthene	15.00	252	548228m	43.60	ng	
88) Benzo(k)fluoranthene	15.02	252	439394m	38.76	ng	
89) Benzo(a)pyrene	15.34	252	458651	41.08	ng	99
90) Dibenzo(a,h)anthracene	16.79	278	403956	38.94	ng	99
91) Benzo(g,h,i)perylene	17.20	276	408735	38.92	ng	# 93

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

