

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
 Data File : BF087655.D
 Acq On : 4 Jun 2016 9:43
 Operator : UM/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF060416

Manual Integrations
 APPROVED

SOHIL
 6/6/2016 7:18:07 PM

Quant Time: Jun 05 01:10:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 04 11:07:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	113462	20.00	ng	-0.01
21) Naphthalene-d8	8.15	136	457228	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	222933	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	398573	20.00	ng	0.00
75) Chrysene-d12	14.02	240	310747	20.00	ng	0.00
86) Perylene-d12	15.44	264	273073	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.45	112	591540	84.27	ng	0.00
7) Phenol-d6	6.49	99	726560	82.53	ng	0.00
23) Nitrobenzene-d5	7.43	82	624812	79.06	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	170666	76.27	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1222392	83.77	ng	0.00
78) Terphenyl-d14	12.97	244	959577	75.35	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.48	88	136755	40.30	ng	# 30
3) Pyridine	3.19	79	384281	42.87	ng	96
4) n-Nitrosodimethylamine	3.14	42	151472	40.00	ng	99
6) Aniline	6.52	93	505207	40.49	ng	99
8) 2-Chlorophenol	6.64	128	310944	38.49	ng	99
9) Benzaldehyde	6.41	77	245409	41.23	ng	98
10) Phenol	6.50	94	453252	45.21	ng	100
11) bis(2-Chloroethyl)ether	6.59	93	273038	37.49	ng	98
12) 1,3-Dichlorobenzene	6.80	146	328032	37.93	ng	99
13) 1,4-Dichlorobenzene	6.88	146	343427	39.57	ng	99
14) 1,2-Dichlorobenzene	7.04	146	314182	38.62	ng	100
15) Benzyl Alcohol	7.00	79	273511	42.05	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.14	45	427336	40.23	ng	99
17) 2-Methylphenol	7.11	107	260684	40.19	ng	99
18) Hexachloroethane	7.38	117	125523	41.36	ng	99
19) n-Nitroso-di-n-propylamine	7.28	70	205238	37.33	ng	99
20) 3+4-Methylphenols	7.27	107	293224	36.90	ng	100
22) Acetophenone	7.28	105	442159	42.70	ng	99
24) Nitrobenzene	7.45	77	344629	43.57	ng	99
25) Isophorone	7.69	82	578199	40.19	ng	99
26) 2-Nitrophenol	7.76	139	144869	39.42	ng	98
27) 2,4-Dimethylphenol	7.80	122	321576	42.22	ng	98
28) bis(2-Chloroethoxy)methane	7.91	93	356563	39.07	ng	100
29) 2,4-Dichlorophenol	8.00	162	249671	39.91	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	269689	41.17	ng	99
31) Naphthalene	8.17	128	898299	40.41	ng	100
32) Benzoic acid	7.92	122	209138	45.52	ng	94
33) 4-Chloroaniline	8.22	127	368966	38.21	ng	100
34) Hexachlorobutadiene	8.30	225	143968	42.27	ng	99
35) Caprolactam	8.59	113	80083	38.97	ng	97
36) 4-Chloro-3-methylphenol	8.70	107	263035	40.65	ng	99
37) 2-Methylnaphthalene	8.87	142	578873	39.43	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	243126	39.57	ng	# 1
40) Hexachlorocyclopentadiene	9.02	237	137672	38.09	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	181690	39.16	nq	100
43) 2,4,5-Trichlorophenol	9.18	196	186739	43.24	nq	99
45) 1,1'-Biphenyl	9.32	154	648419	35.84	nq	99
46) 2-Chloronaphthalene	9.35	162	522615	36.29	nq	99
47) 2-Nitroaniline	9.44	65	150082	37.01	nq	99
48) Acenaphthylene	9.77	152	903053	40.21	nq	99
49) Dimethylphthalate	9.63	163	666679	41.13	nq	100
50) 2,6-Dinitrotoluene	9.69	165	159227	43.18	nq	92
51) Acenaphthene	9.94	154	581879	40.27	nq	100
52) 3-Nitroaniline	9.85	138	155677	36.91	nq	99
53) 2,4-Dinitrophenol	9.95	184	64419	40.51	nq	99
54) Dibenzofuran	10.11	168	810312	41.19	nq	98
55) 4-Nitrophenol	10.00	139	124334	34.49	nq	98
56) 2,4-Dinitrotoluene	10.09	165	215583	45.91	nq	97
57) Fluorene	10.44	166	567212	36.88	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.23	232	144816	38.93	nq #	55
59) Diethylphthalate	10.33	149	624528	38.37	nq	100
60) 4-Chlorophenyl-phenylether	10.44	204	242975	38.84	nq	99
61) 4-Nitroaniline	10.47	138	175516	43.29	nq	98
62) Azobenzene	10.60	77	681516	41.60	nq	99
64) 4,6-Dinitro-2-methylphenol	10.49	198	85445	38.96	nq	99
65) n-Nitrosodiphenylamine	10.56	169	501529	38.42	nq	100
66) 4-Bromophenyl-phenylether	10.94	248	174206	44.08	nq	99
67) Hexachlorobenzene	10.99	284	167233	37.84	nq	99
68) Atrazine	11.10	200	162946	42.14	nq	99
69) Pentachlorophenol	11.19	266	114886	43.75	nq	99
70) Phenanthrene	11.40	178	803427	37.18	nq	100
71) Anthracene	11.46	178	917088	42.35	nq	100
72) Carbazole	11.61	167	773835	37.75	nq	100
73) Di-n-butylphthalate	11.95	149	943583	42.94	nq	100
74) Fluoranthene	12.59	202	889177	41.81	nq	99
76) Benzidine	12.72	184	541619	44.20	nq	99
77) Pyrene	12.82	202	927987	41.94	nq	100
79) Butylbenzylphthalate	13.45	149	422757	44.89	nq	99
80) Benzo(a)anthracene	14.01	228	720385	40.17	nq	99
81) 3,3'-Dichlorobenzidine	13.98	252	230917	45.64	nq	98
82) Chrysene	14.05	228	736276	41.25	nq	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	491745	43.88	nq	99
84) Di-n-octyl phthalate	14.62	149	860068	41.28	nq #	100
85) Indeno(1,2,3-cd)pyrene	16.85	276	637143	43.18	nq #	100
87) Benzo(b)fluoranthene	15.04	252	750116m	42.23	nq	
88) Benzo(k)fluoranthene	15.06	252	556950m	37.68	nq	
89) Benzo(a)pyrene	15.38	252	608505	40.75	nq	99
90) Dibenzo(a,h)anthracene	16.86	278	533660	42.00	nq	99
91) Benzo(g,h,i)perylene	17.26	276	538810	42.03	nq	99

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

