

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062116\
 Data File : BF088233.D
 Acq On : 22 Jun 2016 1:06
 Operator : UM/SJ
 Sample : PB91471BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB91471BS

Manual Integrations
 APPROVED

sohil
 6/22/2016 6:08:02 PM

Quant Time: Jun 22 01:43:05 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:34:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	86251	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	334118	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	157742	20.00	ng	0.00
63) Phenanthrene-d10	11.16	188	290521	20.00	ng	0.00
75) Chrysene-d12	13.79	240	207937	20.00	ng	0.00
86) Perylene-d12	15.17	264	177178	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.26	112	641088	127.07	ng	0.01
7) Phenol-d6	6.31	99	748713	122.45	ng	0.00
23) Nitrobenzene-d5	7.22	82	447746	78.25	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	170235	102.21	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	824795	81.94	ng	0.00
78) Terphenyl-d14	12.75	244	796378	87.15	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.19	88	77820	31.74	ng	# 34
3) Pyridine	2.84	79	221824	38.32	ng	91
4) n-Nitrosodimethylamine	2.80	42	91087	37.16	ng	79
6) Aniline	6.32	93	144834	16.64	ng	# 1
8) 2-Chlorophenol	6.44	128	247956	41.52	ng	82
10) Phenol	6.33	94	245441	38.52	ng	97
11) bis(2-Chloroethyl)ether	6.40	93	238545	44.49	ng	87
12) 1,3-Dichlorobenzene	6.59	146	254308m	39.69	ng	
13) 1,4-Dichlorobenzene	6.67	146	255083	39.52	ng	96
14) 1,2-Dichlorobenzene	6.82	146	248324m	42.31	ng	
15) Benzyl Alcohol	6.81	79	174389	36.46	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.94	45	248773	34.34	ng	# 1
17) 2-Methylphenol	6.94	107	197821	40.85	ng	95
18) Hexachloroethane	7.15	117	90419	39.48	ng	# 82
19) n-Nitroso-di-n-propylamine	7.08	70	153200	39.17	ng	# 77
20) 3+4-Methylphenols	7.08	107	258592	45.76	ng	# 83
22) Acetophenone	7.07	105	331985	44.46	ng	97
24) Nitrobenzene	7.24	77	259267	46.78	ng	96
25) Isophorone	7.48	82	443811	41.88	ng	98
26) 2-Nitrophenol	7.56	139	137874	46.44	ng	# 73
27) 2,4-Dimethylphenol	7.61	122	227507	41.62	ng	93
28) bis(2-Chloroethoxy)methane	7.70	93	304435	46.31	ng	97
29) 2,4-Dichlorophenol	7.80	162	201939	44.24	ng	98
30) 1,2,4-Trichlorobenzene	7.88	180	213014	42.86	ng	98
31) Naphthalene	7.95	128	687824	41.60	ng	99
32) Benzoic acid	7.77	122	117083	38.90	ng	94
33) 4-Chloroaniline	8.02	127	85155	11.53	ng	# 92
34) Hexachlorobutadiene	8.08	225	111416	42.51	ng	100
35) Caprolactam	8.40	113	66212m	43.80	ng	
36) 4-Chloro-3-methylphenol	8.52	107	221957	45.47	ng	91
37) 2-Methylnaphthalene	8.65	142	468721	43.01	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.82	216	195902	49.32	ng	# 1
40) Hexachlorocyclopentadiene	8.80	237	112471	44.20	ng	99
42) 2,4,6-Trichlorophenol	8.94	196	129051	40.91	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.99	196	126785	38.63	nq	97
45) 1,1'-Biphenyl	9.12	154	580700	49.69	nq	96
46) 2-Chloronaphthalene	9.14	162	450176	44.10	nq	94
47) 2-Nitroaniline	9.24	65	126043	41.86	nq	# 82
48) Acenaphthylene	9.55	152	707188	43.56	nq	99
49) Dimethylphthalate	9.43	163	513668	44.95	nq	100
50) 2,6-Dinitrotoluene	9.48	165	112802	43.11	nq	# 62
51) Acenaphthene	9.72	154	418939	41.36	nq	99
52) 3-Nitroaniline	9.66	138	59466	19.69	nq	88
53) 2,4-Dinitrophenol	9.77	184	98974	71.10	nq	98
54) Dibenzofuran	9.90	168	573121	41.09	nq	# 91
55) 4-Nitrophenol	9.85	139	135980	58.02	nq	# 76
56) 2,4-Dinitrotoluene	9.90	165	141760	42.15	nq	# 63
57) Fluorene	10.24	166	478467	51.45	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.02	232	113604	44.05	nq	# 58
59) Diethylphthalate	10.12	149	494950	42.79	nq	99
60) 4-Chlorophenyl-phenylether	10.23	204	186750	40.26	nq	95
61) 4-Nitroaniline	10.27	138	109864	38.36	nq	97
62) Azobenzene	10.39	77	497841	43.53	nq	95
64) 4,6-Dinitro-2-methylphenol	10.31	198	78152	43.31	nq	# 64
65) n-Nitrosodiphenylamine	10.35	169	410167	42.55	nq	99
66) 4-Bromophenyl-phenylether	10.72	248	136609	43.83	nq	97
67) Hexachlorobenzene	10.78	284	133000	41.07	nq	92
68) Atrazine	10.89	200	122552	41.98	nq	98
69) Pentachlorophenol	10.98	266	106503	62.39	nq	97
70) Phenanthrene	11.19	178	659615	43.27	nq	99
71) Anthracene	11.24	178	683393	44.44	nq	99
72) Carbazole	11.40	167	671367	46.65	nq	99
73) Di-n-butylphthalate	11.74	149	773011	42.43	nq	99
74) Fluoranthene	12.38	202	713058	44.32	nq	95
76) Benzidine	12.50	184	208088	36.14	nq	98
77) Pyrene	12.59	202	686862	44.51	nq	99
79) Butylbenzylphthalate	13.22	149	325187	47.67	nq	89
80) Benzo(a)anthracene	13.78	228	493415	41.64	nq	99
81) 3,3'-Dichlorobenzidine	13.75	252	108415	34.02	nq	# 94
82) Chrysene	13.82	228	550918	49.42	nq	98
83) Bis(2-ethylhexyl)phthalate	13.78	149	383474	46.18	nq	99
84) Di-n-octyl phthalate	14.39	149	703107	52.10	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.46	276	463111	44.71	nq	# 100
87) Benzo(b)fluoranthene	14.79	252	449456m	36.40	nq	
88) Benzo(k)fluoranthene	14.82	252	477798m	51.29	nq	
89) Benzo(a)pyrene	15.11	252	445323	44.57	nq	99
90) Dibenzo(a,h)anthracene	16.46	278	383796	41.36	nq	99
91) Benzo(g,h,i)perylene	16.83	276	406711	42.61	nq	98

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

