

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022417\
 Data File : BF093180.D
 Acq On : 25 Feb 2017 9:20
 Operator : SJ/MA
 Sample : I1972-21MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Z5-BASEMSD

Manual Integrations
 APPROVED

mohammad
 2/27/2017 1:08:29 PM

Quant Time: Feb 27 00:47:41 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 17:31:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	153131	20.00	ng	-0.02
21) Naphthalene-d8	8.12	136	588875	20.00	ng	-0.02
38) Acenaphthene-d10	9.88	164	259018	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	386160	20.00	ng	-0.01
75) Chrysene-d12	14.01	240	317100	20.00	ng	-0.01
86) Perylene-d12	15.44	264	284486	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1041046	116.63	ng	-0.01
7) Phenol-d6	6.49	99	1371932	119.46	ng	0.00
23) Nitrobenzene-d5	7.41	82	920992	87.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	179123	95.61	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	1328349	92.91	ng	-0.01
78) Terphenyl-d14	12.96	244	762322	56.49	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.44	88	174736	36.23	ng	# 85
3) Pyridine	3.16	79	363788	29.66	ng	88
4) n-Nitrosodimethylamine	3.12	42	247233	42.93	ng	88
6) Aniline	6.50	93	192138	12.26	ng	# 12
8) 2-Chlorophenol	6.62	128	409845	41.62	ng	94
9) Benzaldehyde	6.37	77	177840	21.61	ng	86
10) Phenol	6.50	94	652308	48.55	ng	# 77
11) bis(2-Chloroethyl)ether	6.58	93	463161	43.19	ng	97
12) 1,3-Dichlorobenzene	6.77	146	465132	40.33	ng	# 94
13) 1,4-Dichlorobenzene	6.85	146	483978	41.46	ng	96
14) 1,2-Dichlorobenzene	7.00	146	414762	37.16	ng	97
15) Benzyl Alcohol	6.99	79	353043	41.52	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.12	45	743701	39.91	ng	93
17) 2-Methylphenol	7.10	107	296102	35.05	ng	98
18) Hexachloroethane	7.34	117	141401	35.48	ng	94
19) n-Nitroso-di-n-propylamine	7.25	70	309653	42.36	ng	93
20) 3+4-Methylphenols	7.26	107	400531	39.66	ng	# 75
22) Acetophenone	7.25	105	614701	45.72	ng	# 97
24) Nitrobenzene	7.42	77	450000	41.44	ng	99
25) Isophorone	7.66	82	849974	43.13	ng	97
26) 2-Nitrophenol	7.74	139	228823	44.33	ng	89
27) 2,4-Dimethylphenol	7.79	122	284229	31.36	ng	92
28) bis(2-Chloroethoxy)methane	7.88	93	569341	43.63	ng	100
29) 2,4-Dichlorophenol	7.98	162	335905	39.73	ng	85
30) 1,2,4-Trichlorobenzene	8.06	180	358339	39.33	ng	95
31) Naphthalene	8.14	128	1156786	38.75	ng	99
32) Benzoic acid	7.93	122	1281m	7.38	ng	
33) 4-Chloroaniline	8.20	127	125525	10.72	ng	99
34) Hexachlorobutadiene	8.26	225	183518	36.61	ng	98
35) Caprolactam	8.57	113	110645	41.61	ng	# 36
36) 4-Chloro-3-methylphenol	8.69	107	337203	38.26	ng	97
37) 2-Methylnaphthalene	8.84	142	774374m	40.40	ng	
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	315932	43.45	ng	99
40) Hexachlorocyclopentadiene	8.99	237	76538	23.33	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	215885	45.19	nq	94
43) 2,4,5-Trichlorophenol	9.17	196	220865	42.19	nq	# 87
45) 1,1'-Biphenyl	9.30	154	817876	43.61	nq	96
46) 2-Chloronaphthalene	9.32	162	668543	45.57	nq	94
47) 2-Nitroaniline	9.42	65	216459	43.88	nq	88
48) Acenaphthylene	9.74	152	1065743	42.05	nq	99
49) Dimethylphthalate	9.61	163	1030214	59.05	nq	99
50) 2,6-Dinitrotoluene	9.68	165	185300	49.18	nq	93
51) Acenaphthene	9.92	154	626719	41.36	nq	97
52) 3-Nitroaniline	9.84	138	127265	29.51	nq	86
53) 2,4-Dinitrophenol	9.96	184	5310	7.34	nq	# 64
54) Dibenzofuran	10.09	168	906014	44.70	nq	93
55) 4-Nitrophenol	10.02	139	184336	59.36	nq	87
56) 2,4-Dinitrotoluene	10.08	165	183930	36.67	nq	# 87
57) Fluorene	10.43	166	685971	43.99	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.21	232	148281	42.46	nq	94
59) Diethylphthalate	10.30	149	737795	44.87	nq	99
60) 4-Chlorophenyl-phenylether	10.42	204	291596	38.92	nq	92
61) 4-Nitroaniline	10.45	138	137426	31.12	nq	93
62) Azobenzene	10.58	77	776136	45.69	nq	97
64) 4,6-Dinitro-2-methylphenol	10.49	198	24316	12.44	nq	# 29
65) n-Nitrosodiphenylamine	10.54	169	602775	48.86	nq	100
66) 4-Bromophenyl-phenylether	10.91	248	178925	44.35	nq	94
67) Hexachlorobenzene	10.98	284	162366	40.73	nq	96
68) Atrazine	11.07	200	162458	41.04	nq	100
69) Pentachlorophenol	11.17	266	121683	61.00	nq	97
70) Phenanthrene	11.39	178	879645	39.76	nq	98
71) Anthracene	11.45	178	840443	37.83	nq	97
72) Carbazole	11.60	167	738570	34.78	nq	99
73) Di-n-butylphthalate	11.93	149	1007434	43.86	nq	# 98
74) Fluoranthene	12.58	202	813129	35.63	nq	97
76) Benzidine	12.71	184	144671	10.71	nq	97
77) Pyrene	12.81	202	796434	33.56	nq	99
79) Butylbenzylphthalate	13.43	149	366794	38.06	nq	96
80) Benzo(a)anthracene	14.00	228	711011	36.66	nq	99
81) 3,3'-Dichlorobenzidine	13.96	252	176421	25.52	nq	# 96
82) Chrysene	14.03	228	600336	33.06	nq	100
83) Bis(2-ethylhexyl)phthalate	13.99	149	510802	38.76	nq	# 97
84) Di-n-octyl phthalate	14.59	149	895816	41.74	nq	98
85) Indeno(1,2,3-cd)pyrene	16.85	276	534472	38.00	nq	# 92
87) Benzo(b)fluoranthene	15.04	252	814836m	43.52	nq	
88) Benzo(k)fluoranthene	15.06	252	506349m	31.32	nq	
89) Benzo(a)pyrene	15.38	252	589904	36.42	nq	# 96
90) Dibenzo(a,h)anthracene	16.87	278	467423	35.71	nq	98
91) Benzo(g,h,i)perylene	17.28	276	421099	31.84	nq	# 94

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

