

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
 Data File : BF086729.D
 Acq On : 6 May 2016 17:40
 Operator : UM/SJ
 Sample : PB90268BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB90268BS

Manual Integrations
 APPROVED

Sohil
 5/9/2016 7:17:12 PM

Quant Time: May 07 03:36:08 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 07 03:18:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	154577	40.00	ng	0.00
21) Naphthalene-d8	8.00	136	640054	40.00	ng	0.00
38) Acenaphthene-d10	9.74	164	300091	40.00	ng	0.00
63) Phenanthrene-d10	11.23	188	562973	40.00	ng	0.00
75) Chrysene-d12	13.86	240	421247	40.00	ng	0.00
86) Perylene-d12	15.23	264	316861	40.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	1065105	114.07	ng	0.01
7) Phenol-d6	6.38	99	1502905	119.49	ng	0.00
23) Nitrobenzene-d5	7.29	82	1014944	81.49	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	427429	145.60	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1702503	93.00	ng	0.00
78) Terphenyl-d14	12.81	244	1346221	64.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.35	88	156140	32.35	ng	# 75
3) Pyridine	3.16	79	391263m	33.06	ng	
4) n-Nitrosodimethylamine	2.99	42	292043	40.69	ng	# 52
6) Aniline	6.38	93	223687	14.41	ng	# 1
8) 2-Chlorophenol	6.50	128	424052	37.84	ng	# 79
9) Benzaldehyde	6.27	77	47921	7.64	ng	96
10) Phenol	6.40	94	575003	41.79	ng	84
11) bis(2-Chloroethyl)ether	6.45	93	415672	35.10	ng	84
12) 1,3-Dichlorobenzene	6.65	146	459259m	38.60	ng	
13) 1,4-Dichlorobenzene	6.73	146	473051	38.45	ng	97
14) 1,2-Dichlorobenzene	6.88	146	389758	36.39	ng	96
15) Benzyl Alcohol	6.88	79	370172	45.49	ng	89
16) 2,2'-oxybis(1-Chloropropan	7.00	45	853430	35.24	ng	81
17) 2-Methylphenol	6.99	107	343789	38.11	ng	# 80
18) Hexachloroethane	7.22	117	178958	38.53	ng	93
19) n-Nitroso-di-n-propylamine	7.15	70	343664	39.99	ng	# 78
20) 3+4-Methylphenols	7.15	107	443290	41.96	ng	# 67
22) Acetophenone	7.14	105	613915	41.16	ng	99
24) Nitrobenzene	7.30	77	455318	35.36	ng	95
25) Isophorone	7.56	82	901830	38.88	ng	97
26) 2-Nitrophenol	7.62	139	228829	39.83	ng	# 75
27) 2,4-Dimethylphenol	7.68	122	413548	40.99	ng	92
28) bis(2-Chloroethoxy)methane	7.76	93	522273	40.91	ng	98
29) 2,4-Dichlorophenol	7.87	162	378806	43.67	ng	99
30) 1,2,4-Trichlorobenzene	7.94	180	384979	39.78	ng	97
31) Naphthalene	8.02	128	1302163	39.82	ng	99
32) Benzoic acid	7.84	122	263901	63.71	ng	# 85
33) 4-Chloroaniline	8.09	127	85895	6.78	ng	96
34) Hexachlorobutadiene	8.13	225	209883	38.20	ng	99
35) Caprolactam	8.50	113	122473m	45.61	ng	
36) 4-Chloro-3-methylphenol	8.58	107	396817	40.10	ng	92
37) 2-Methylnaphthalene	8.70	142	796223	41.27	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.88	216	357154	41.12	ng	# 97
40) Hexachlorocyclopentadiene	8.86	237	452259	110.43	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.00	196	263299	47.67	ng	92
43) 2,4,5-Trichlorophenol	9.05	196	256404	44.43	ng #	90
45) 1,1'-Biphenyl	9.17	154	999229	40.73	ng	97
46) 2-Chloronaphthalene	9.20	162	746668	39.44	ng	94
47) 2-Nitroaniline	9.31	65	280811	41.72	ng	97
48) Acenaphthylene	9.61	152	1092052	38.86	ng	99
49) Dimethylphthalate	9.49	163	960211	42.92	ng	100
50) 2,6-Dinitrotoluene	9.55	165	221423	46.12	ng	90
51) Acenaphthene	9.78	154	689681	37.20	ng	99
52) 3-Nitroaniline	9.72	138	76100	13.94	ng	97
53) 2,4-Dinitrophenol	9.82	184	220369	97.72	ng #	76
54) Dibenzofuran	9.95	168	978508	42.58	ng	92
55) 4-Nitrophenol	9.90	139	245183	77.05	ng #	40
56) 2,4-Dinitrotoluene	9.95	165	241428	40.58	ng #	71
57) Fluorene	10.29	166	720564	38.69	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.08	232	205400	48.46	ng	96
59) Diethylphthalate	10.19	149	938948	44.34	ng	97
60) 4-Chlorophenyl-phenylether	10.29	204	384036	43.37	ng #	85
61) 4-Nitroaniline	10.34	138	177813	33.91	ng #	70
62) Azobenzene	10.45	77	1065442	45.41	ng	88
64) 4,6-Dinitro-2-methylphenol	10.36	198	164993	55.08	ng	92
65) n-Nitrosodiphenylamine	10.42	169	738659	41.32	ng	99
66) 4-Bromophenyl-phenylether	10.77	248	215768	38.57	ng #	86
67) Hexachlorobenzene	10.84	284	284529	41.70	ng	99
68) Atrazine	10.96	200	247208	45.13	ng	95
69) Pentachlorophenol	11.05	266	315419	94.16	ng	97
70) Phenanthrene	11.25	178	1260022	41.67	ng	99
71) Anthracene	11.30	178	1180905	37.29	ng	100
72) Carbazole	11.47	167	1127194	40.99	ng	99
73) Di-n-butylphthalate	11.80	149	1355868	41.53	ng #	99
74) Fluoranthene	12.43	202	1153645	38.07	ng	99
76) Benzidine	12.57	184	313088	30.22	ng	96
77) Pyrene	12.66	202	1146348	34.13	ng	99
79) Butylbenzylphthalate	13.29	149	510586	36.30	ng #	69
80) Benzo(a)anthracene	13.85	228	927933	38.95	ng	99
81) 3,3'-Dichlorobenzidine	13.81	252	115616	7.72	ng	99
82) Chrysene	13.88	228	902955	35.53	ng	98
83) Bis(2-ethylhexyl)phthalate	13.85	149	638247	39.21	ng #	95
84) Di-n-octyl phthalate	14.45	149	1125983	47.01	ng #	89
85) Indeno(1,2,3-cd)pyrene	16.55	276	692191	31.74	ng	95
87) Benzo(b)fluoranthene	14.85	252	797810m	41.90	ng	
88) Benzo(k)fluoranthene	14.88	252	803898m	44.81	ng	
89) Benzo(a)pyrene	15.19	252	749258	42.72	ng	97
90) Dibenzo(a,h)anthracene	16.56	278	580019	34.43	ng #	93
91) Benzo(g,h,i)perylene	16.93	276	589322	33.32	ng #	90

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

