

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020918\
 Data File : BF102893.D
 Acq On : 9 Feb 2018 17:16
 Operator : SJ/JU
 Sample : PB106115BSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB106115BSD

Manual Integrations
 APPROVED

Sohil
 2/13/2018 4:36:49 AM

Quant Time: Feb 09 22:21:29 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	162584	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	668410	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	280868	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	405544	20.00	ng	0.00
75) Chrysene-d12	14.06	240	320969	20.00	ng	0.01
86) Perylene-d12	15.54	264	311962	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	1214390	113.43	ng	0.02
7) Phenol-d6	6.52	99	1453912	112.79	ng	0.01
23) Nitrobenzene-d5	7.45	82	960685	99.70	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	243787	107.84	ng	0.01
44) 2-Fluorobiphenyl	9.25	172	1441994	85.19	ng	0.00
78) Terphenyl-d14	13.00	244	1095317	80.80	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.72	88	197619	37.373	ng	87
3) Pyridine	3.50	79	590206	40.399	ng	90
4) n-Nitrosodimethylamine	3.45	42	300403	48.059	ng	92
6) Aniline	6.55	93	342289	17.981	ng	95
8) 2-Chlorophenol	6.67	128	449767	40.808	ng	96
9) Benzaldehyde	6.44	77	231052	24.136	ng	98
10) Phenol	6.53	94	580132	41.995	ng	98
11) bis(2-Chloroethyl)ether	6.62	93	468407	40.748	ng	92
12) 1,3-Dichlorobenzene	6.83	146	478206	37.467	ng	97
13) 1,4-Dichlorobenzene	6.90	146	482808	37.975	ng	99
14) 1,2-Dichlorobenzene	7.06	146	447985	37.830	ng	100
15) Benzyl Alcohol	7.03	79	418628	42.241	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.16	45	820881	37.592	ng	99
17) 2-Methylphenol	7.14	107	388413	38.205	ng	97
18) Hexachloroethane	7.40	117	168576	38.666	ng	94
19) n-Nitroso-di-n-propylamine	7.30	70	322016	37.889	ng	93
20) 3+4-Methylphenols	7.29	107	456294	36.396	ng	# 83
22) Acetophenone	7.30	105	618206	37.795	ng	# 93
24) Nitrobenzene	7.47	77	513498	45.307	ng	95
25) Isophorone	7.70	82	957334	43.149	ng	95
26) 2-Nitrophenol	7.79	139	234231	49.888	ng	98
27) 2,4-Dimethylphenol	7.82	122	426619	45.513	ng	98
28) bis(2-Chloroethoxy)methane	7.92	93	579732	44.109	ng	100
29) 2,4-Dichlorophenol	8.03	162	360675	42.327	ng	98
30) 1,2,4-Trichlorobenzene	8.11	180	362887	37.645	ng	100
31) Naphthalene	8.19	128	1246619	38.173	ng	100
32) Benzoic acid	7.93	122	247433	40.212	ng	94
33) 4-Chloroaniline	8.23	127	137201	9.752	ng	95
34) Hexachlorobutadiene	8.30	225	185165	37.485	ng	98
35) Caprolactam	8.61	113	125579m	42.436	ng	
36) 4-Chloro-3-methylphenol	8.72	107	402358	42.025	ng	99
37) 2-Methylnaphthalene	8.89	142	813750	38.059	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.05	216	304118	39.767	ng	97
40) Hexachlorocyclopentadiene	9.03	237	88182	24.256	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.16	196	218384	43.476	ng	99
43) 2,4,5-Trichlorophenol	9.20	196	230272	40.673	ng	98
45) 1,1'-Biphenyl	9.35	154	920869	38.926	ng	99
46) 2-Chloronaphthalene	9.37	162	719837	38.651	ng	99
47) 2-Nitroaniline	9.47	65	287599	49.895	ng	86
48) Acenaphthylene	9.79	152	1075366	37.176	ng	99
49) Dimethylphthalate	9.65	163	885495	40.434	ng	100
50) 2,6-Dinitrotoluene	9.71	165	182767	43.866	ng	90
51) Acenaphthene	9.96	154	641866	37.104	ng	99
52) 3-Nitroaniline	9.87	138	114211	22.278	ng	92
53) 2,4-Dinitrophenol	9.99	184	34409	39.692	ng	# 18
54) Dibenzofuran	10.13	168	929222	38.116	ng	96
55) 4-Nitrophenol	10.04	139	288219	68.995	ng	91
56) 2,4-Dinitrotoluene	10.12	165	229771	41.560	ng	97
57) Fluorene	10.47	166	678949	38.278	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.25	232	155860	39.720	ng	95
59) Diethylphthalate	10.35	149	853849	39.486	ng	99
60) 4-Chlorophenyl-phenylether	10.46	204	313807	39.993	ng	95
61) 4-Nitroaniline	10.49	138	169763	34.789	ng	99
62) Azobenzene	10.63	77	870156	40.280	ng	96
64) 4,6-Dinitro-2-methylphenol	10.52	198	30057	24.459	ng	91
65) n-Nitrosodiphenylamine	10.59	169	621294	43.433	ng	99
66) 4-Bromophenyl-phenylether	10.96	248	184917	43.105	ng	94
67) Hexachlorobenzene	11.03	284	180298	38.097	ng	93
68) Atrazine	11.11	200	181936	43.112	ng	95
69) Pentachlorophenol	11.22	266	175093	63.026	ng	100
70) Phenanthrene	11.44	178	866239	38.353	ng	99
71) Anthracene	11.49	178	898422	39.109	ng	99
72) Carbazole	11.65	167	843639	37.486	ng	99
73) Di-n-butylphthalate	11.97	149	1191487	43.583	ng	100
74) Fluoranthene	12.63	202	870844	38.322	ng	99
76) Benzidine	12.75	184	513848	34.833	ng	98
77) Pyrene	12.86	202	900481	35.777	ng	99
79) Butylbenzylphthalate	13.47	149	504790	42.058	ng	94
80) Benzo(a)anthracene	14.05	228	825896	40.914	ng	99
81) 3,3'-Dichlorobenzidine	14.01	252	297515	39.751	ng	96
82) Chrysene	14.09	228	776490	38.160	ng	100
83) Bis(2-ethylhexyl)phthalate	14.03	149	727062	47.156	ng	100
84) Di-n-octyl phthalate	14.64	149	1265256	54.653	ng	97
85) Indeno(1,2,3-cd)pyrene	17.06	276	653410	38.717	ng	98
87) Benzo(b)fluoranthene	15.11	252	808570	39.977	ng	98
88) Benzo(k)fluoranthene	15.14	252	772822	39.990	ng	100
89) Benzo(a)pyrene	15.49	252	744601	40.900	ng	98
90) Dibenzo(a,h)anthracene	17.07	278	552628	37.398	ng	98
91) Benzo(g,h,i)perylene	17.51	276	513570	35.508	ng	97

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

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