

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085554.D
 Acq On : 19 Mar 2016 5:12
 Operator : UM/SJ
 Sample : PB89025BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB89025BS

Manual Integrations
 APPROVED

apatel
 3/21/2016 5:08:06 PM

Quant Time: Mar 19 06:37:11 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	127009	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	515176	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	242112	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	436226	20.00	ng	-0.01
75) Chrysene-d12	14.01	240	385238	20.00	ng	0.00
86) Perylene-d12	15.43	264	282626	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	849895	112.40	ng	0.01
7) Phenol-d6	6.49	99	1129459	116.46	ng	0.00
23) Nitrobenzene-d5	7.42	82	671136	75.62	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	297061	125.26	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1198435	84.25	ng	-0.01
78) Terphenyl-d14	12.96	244	1208464	75.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.55	88	124316	33.82	ng	99
3) Pyridine	3.28	79	301544	33.41	ng	100
4) n-Nitrosodimethylamine	3.22	42	138509	36.76	ng	95
6) Aniline	6.52	93	304217	23.36	ng	98
8) 2-Chlorophenol	6.64	128	362039	41.39	ng	97
9) Benzaldehyde	6.40	77	52519	8.61	ng	86
10) Phenol	6.50	94	510671	46.23	ng	95
11) bis(2-Chloroethyl)ether	6.60	93	359684	43.19	ng	95
12) 1,3-Dichlorobenzene	6.79	146	354358m	37.17	ng	
13) 1,4-Dichlorobenzene	6.87	146	376191	38.69	ng	98
14) 1,2-Dichlorobenzene	7.03	146	346474	40.13	ng	96
15) Benzyl Alcohol	7.00	79	272368	41.00	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	416449	38.03	ng	99
17) 2-Methylphenol	7.11	107	293140	40.43	ng	99
18) Hexachloroethane	7.36	117	136708	38.94	ng	# 81
19) n-Nitroso-di-n-propylamine	7.27	70	212597	36.70	ng	92
20) 3+4-Methylphenols	7.27	107	362884	43.21	ng	89
22) Acetophenone	7.27	105	454263	37.16	ng	# 96
24) Nitrobenzene	7.44	77	404100	41.59	ng	91
25) Isophorone	7.68	82	698394	39.30	ng	96
26) 2-Nitrophenol	7.76	139	193961	40.40	ng	# 86
27) 2,4-Dimethylphenol	7.80	122	342901	40.39	ng	97
28) bis(2-Chloroethoxy)methane	7.89	93	415290	41.09	ng	98
29) 2,4-Dichlorophenol	8.00	162	300645	42.88	ng	95
30) 1,2,4-Trichlorobenzene	8.08	180	303393	38.63	ng	96
31) Naphthalene	8.16	128	1019045	39.19	ng	100
32) Benzoic acid	7.92	122	229094	32.27	ng	99
33) 4-Chloroaniline	8.21	127	117227	11.00	ng	# 93
34) Hexachlorobutadiene	8.28	225	157346	37.59	ng	99
35) Caprolactam	8.58	113	99191m	41.12	ng	
36) 4-Chloro-3-methylphenol	8.70	107	332781	40.64	ng	98
37) 2-Methylnaphthalene	8.86	142	693252	43.74	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	265290	36.12	ng	98
40) Hexachlorocyclopentadiene	9.01	237	297996	90.03	ng	100

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42) 2,4,6-Trichlorophenol	9.13	196	210879	41.95	ng	100
43) 2,4,5-Trichlorophenol	9.18	196	218984	42.80	ng #	94
45) 1,1'-Biphenyl	9.32	154	734975	35.19	ng	98
46) 2-Chloronaphthalene	9.34	162	566554	37.14	ng #	92
47) 2-Nitroaniline	9.44	65	203610	44.12	ng #	84
48) Acenaphthylene	9.76	152	1019964	41.31	ng	99
49) Dimethylphthalate	9.62	163	770040	42.15	ng	99
50) 2,6-Dinitrotoluene	9.68	165	163448	42.38	ng	88
51) Acenaphthene	9.93	154	706369	45.59	ng	99
52) 3-Nitroaniline	9.85	138	90844	18.81	ng #	76
53) 2,4-Dinitrophenol	9.96	184	178899	75.71	ng	95
54) Dibenzofuran	10.10	168	898077	47.60	ng	98
55) 4-Nitrophenol	10.01	139	267626	71.68	ng	93
56) 2,4-Dinitrotoluene	10.08	165	208815	41.35	ng #	49
57) Fluorene	10.45	166	671059	42.68	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.22	232	169646	46.05	ng	98
59) Diethylphthalate	10.32	149	735467	40.49	ng	99
60) 4-Chlorophenyl-phenylether	10.44	204	317451	41.34	ng	92
61) 4-Nitroaniline	10.47	138	200376	42.08	ng	94
62) Azobenzene	10.60	77	794589	45.59	ng	97
64) 4,6-Dinitro-2-methylphenol	10.49	198	117150	41.97	ng	86
65) n-Nitrosodiphenylamine	10.55	169	583464	42.94	ng	99
66) 4-Bromophenyl-phenylether	10.93	248	195106	44.77	ng #	91
67) Hexachlorobenzene	11.00	284	202822	43.81	ng #	88
68) Atrazine	11.09	200	197912	47.84	ng	96
69) Pentachlorophenol	11.19	266	243779	86.73	ng	99
70) Phenanthrene	11.41	178	1054616	45.85	ng	99
71) Anthracene	11.45	178	952837	39.51	ng	100
72) Carbazole	11.61	167	944821	45.41	ng	98
73) Di-n-butylphthalate	11.94	149	1165782	44.74	ng	99
74) Fluoranthene	12.58	202	955609	39.68	ng	97
76) Benzidine	12.71	184	420235	32.44	ng	98
77) Pyrene	12.81	202	963353	34.83	ng	99
79) Butylbenzylphthalate	13.43	149	467457	35.45	ng #	75
80) Benzo(a)anthracene	14.00	228	878103	39.26	ng	99
81) 3,3'-Dichlorobenzidine	13.97	252	130329	15.90	ng #	95
82) Chrysene	14.04	228	685746	31.87	ng	99
83) Bis(2-ethylhexyl)phthalate	13.99	149	582533	35.56	ng #	98
84) Di-n-octyl phthalate	14.61	149	1039606	38.95	ng	98
85) Indeno(1,2,3-cd)pyrene	16.84	276	608012	33.82	ng	99
87) Benzo(b)fluoranthene	15.03	252	777993	42.19	ng #	97
88) Benzo(k)fluoranthene	15.05	252	659524m	43.41	ng	
89) Benzo(a)pyrene	15.37	252	670540	42.89	ng	97
90) Dibenzo(a,h)anthracene	16.85	278	513779	39.40	ng	98
91) Benzo(g,h,i)perylene	17.26	276	494148	39.15	ng #	94

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

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