

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089453.D
 Acq On : 30 Jul 2016 15:14
 Operator : UM/SJ
 Sample : PB92568BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92568BS

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:18:18 PM

Quant Time: Aug 01 04:50:02 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 04:11:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	48927	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	200231	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	91481	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	179734	20.00	ng	0.00
75) Chrysene-d12	13.62	240	100438	20.00	ng	0.00
86) Perylene-d12	14.95	264	83938	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.08	112	405934	132.48	ng	0.02
7) Phenol-d6	6.17	99	489126	120.79	ng	0.00
23) Nitrobenzene-d5	7.08	82	331897	97.84	ng	0.00
41) 2,4,6-Tribromophenol	10.33	330	104395	137.79	ng	0.01
44) 2-Fluorobiphenyl	8.87	172	581366	93.26	ng	0.00
78) Terphenyl-d14	12.58	244	437478	101.94	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.98	88	47731	33.84	ng	# 91
3) Pyridine	2.56	79	121050	40.38	ng	98
4) n-Nitrosodimethylamine	2.54	42	55836	45.68	ng	99
6) Aniline	6.17	93	130726	29.90	ng	# 84
8) 2-Chlorophenol	6.28	128	154049	44.44	ng	83
9) Benzaldehyde	6.04	77	70190	35.50	ng	99
10) Phenol	6.18	94	197347	44.16	ng	94
11) bis(2-Chloroethyl)ether	6.25	93	154625	40.73	ng	91
12) 1,3-Dichlorobenzene	6.44	146	148303m	40.90	ng	
13) 1,4-Dichlorobenzene	6.52	146	161202	40.29	ng	93
14) 1,2-Dichlorobenzene	6.67	146	162453	43.34	ng	97
15) Benzyl Alcohol	6.66	79	129077	51.36	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.80	45	170276	41.12	ng	97
17) 2-Methylphenol	6.79	107	119995	44.79	ng	97
18) Hexachloroethane	7.00	117	62089	40.47	ng	95
19) n-Nitroso-di-n-propylamine	6.94	70	112302	45.99	ng	# 90
20) 3+4-Methylphenols	6.95	107	163563	47.54	ng	# 89
22) Acetophenone	6.92	105	204647	42.93	ng	# 96
24) Nitrobenzene	7.10	77	161882	41.93	ng	91
25) Isophorone	7.34	82	315297	46.14	ng	97
26) 2-Nitrophenol	7.42	139	86080	49.87	ng	94
27) 2,4-Dimethylphenol	7.47	122	141931	51.61	ng	98
28) bis(2-Chloroethoxy)methane	7.56	93	204436	54.05	ng	98
29) 2,4-Dichlorophenol	7.67	162	128927	51.45	ng	96
30) 1,2,4-Trichlorobenzene	7.74	180	135905	47.54	ng	98
31) Naphthalene	7.82	128	459985	46.39	ng	99
32) Benzoic acid	7.62	122	55472	28.95	ng	98
33) 4-Chloroaniline	7.88	127	48748	12.91	ng	96
34) Hexachlorobutadiene	7.93	225	67902	41.47	ng	98
35) Caprolactam	8.26	113	38206m	50.52	ng	
36) 4-Chloro-3-methylphenol	8.38	107	146994	49.36	ng	99
37) 2-Methylnaphthalene	8.50	142	298117	49.90	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.67	216	123132	38.46	ng	99
40) Hexachlorocyclopentadiene	8.65	237	89067	95.64	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.79	196	82517	54.21	nq	98
43) 2,4,5-Trichlorophenol	8.83	196	93449	48.47	nq	# 86
45) 1,1'-Biphenyl	8.97	154	345426	44.35	nq	95
46) 2-Chloronaphthalene	8.99	162	279367	47.76	nq	91
47) 2-Nitroaniline	9.10	65	90940	52.50	nq	# 83
48) Acenaphthylene	9.39	152	432489	46.94	nq	99
49) Dimethylphthalate	9.28	163	328991	47.25	nq	98
50) 2,6-Dinitrotoluene	9.35	165	71325	49.86	nq	# 79
51) Acenaphthene	9.56	154	248316	46.15	nq	99
52) 3-Nitroaniline	9.51	138	33455	22.16	nq	85
53) 2,4-Dinitrophenol	9.63	184	45341	72.23	nq	# 82
54) Dibenzofuran	9.74	168	380408	51.88	nq	96
55) 4-Nitrophenol	9.70	139	83359m	110.80	nq	
56) 2,4-Dinitrotoluene	9.75	165	98319	52.21	nq	# 82
57) Fluorene	10.08	166	303435	47.41	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.87	232	71870	55.83	nq	98
59) Diethylphthalate	9.98	149	337422	47.80	nq	99
60) 4-Chlorophenyl-phenylether	10.08	204	144313	49.16	nq	# 88
61) 4-Nitroaniline	10.12	138	67949	43.67	nq	94
62) Azobenzene	10.24	77	396151	54.50	nq	90
64) 4,6-Dinitro-2-methylphenol	10.16	198	43421	40.75	nq	96
65) n-Nitrosodiphenylamine	10.20	169	284926	50.81	nq	99
66) 4-Bromophenyl-phenylether	10.56	248	82067	47.82	nq	# 81
67) Hexachlorobenzene	10.63	284	84328	47.95	nq	# 71
68) Atrazine	10.74	200	85709	49.75	nq	97
69) Pentachlorophenol	10.82	266	71122	83.21	nq	99
70) Phenanthrene	11.03	178	422371	45.56	nq	99
71) Anthracene	11.08	178	480128	46.59	nq	99
72) Carbazole	11.24	167	413763	47.69	nq	98
73) Di-n-butylphthalate	11.59	149	544766	46.80	nq	99
74) Fluoranthene	12.20	202	430867	44.19	nq	96
76) Benzidine	12.34	184	125255	33.75	nq	97
77) Pyrene	12.43	202	423262	55.60	nq	100
79) Butylbenzylphthalate	13.06	149	173504	50.13	nq	92
80) Benzo(a)anthracene	13.61	228	275439	48.59	nq	99
81) 3,3'-Dichlorobenzidine	13.59	252	55425	27.17	nq	99
82) Chrysene	13.65	228	263664	43.70	nq	100
83) Bis(2-ethylhexyl)phthalate	13.63	149	227217	45.42	nq	# 94
84) Di-n-octyl phthalate	14.24	149	337332	43.21	nq	99
85) Indeno(1,2,3-cd)pyrene	16.14	276	270195	62.99	nq	98
87) Benzo(b)fluoranthene	14.61	252	209837m	40.25	nq	
88) Benzo(k)fluoranthene	14.63	252	228343m	47.65	nq	
89) Benzo(a)pyrene	14.90	252	224894	48.09	nq	98
90) Dibenzo(a,h)anthracene	16.14	278	223418	60.65	nq	100
91) Benzo(g,h,i)perylene	16.48	276	233632	64.52	nq	99

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

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