

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080916\
 Data File : BF089698.D
 Acq On : 9 Aug 2016 20:30
 Operator : UM/SJ
 Sample : PB92711BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92711BS

Manual Integrations
 APPROVED

sohil
 8/10/2016 9:23:10 AM

Quant Time: Aug 10 02:08:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 01:39:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	46259	20.00	ng	-0.03
21) Naphthalene-d8	9.45	136	199366	20.00	ng	-0.02
38) Acenaphthene-d10	12.26	164	90423	20.00	ng	-0.03
63) Phenanthrene-d10	14.66	188	167052	20.00	ng	-0.03
75) Chrysene-d12	18.37	240	101826	20.00	ng	-0.02
86) Perylene-d12	20.02	264	95421	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	337252	122.19	ng	0.02
7) Phenol-d6	6.99	99	439340	117.34	ng	0.01
23) Nitrobenzene-d5	8.33	82	282525	78.38	ng	-0.03
41) 2,4,6-Tribromophenol	13.57	330	81273	108.91	ng	-0.02
44) 2-Fluorobiphenyl	11.22	172	519548	74.67	ng	-0.03
78) Terphenyl-d14	17.03	244	402384	78.01	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	1.80	88	44864	28.36	ng	# 77
3) Pyridine	2.36	79	114731	39.46	ng	98
4) n-Nitrosodimethylamine	2.33	42	49106	39.83	ng	100
6) Aniline	6.91	93	82499	17.00	ng	99
8) 2-Chlorophenol	7.11	128	138056	40.74	ng	99
9) Benzaldehyde	6.72	77	44594m	32.83	ng	
10) Phenol	7.00	94	163119	42.44	ng	98
11) bis(2-Chloroethyl)ether	7.06	93	134156	40.62	ng	99
12) 1,3-Dichlorobenzene	7.31	146	134083	36.43	ng	96
13) 1,4-Dichlorobenzene	7.44	146	137444	37.45	ng	98
14) 1,2-Dichlorobenzene	7.68	146	141949	36.69	ng	96
15) Benzyl Alcohol	7.71	79	109493	44.60	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.92	45	154843	36.82	ng	# 52
17) 2-Methylphenol	7.95	107	108663	44.40	ng	# 87
18) Hexachloroethane	8.19	117	55563	35.92	ng	95
19) n-Nitroso-di-n-propylamine	8.15	70	105001	38.82	ng	96
20) 3+4-Methylphenols	8.22	107	137147	44.40	ng	96
22) Acetophenone	8.10	105	165656	34.85	ng	99
24) Nitrobenzene	8.36	77	139596	40.78	ng	91
25) Isophorone	8.76	82	275519	39.92	ng	98
26) 2-Nitrophenol	8.87	139	64419	41.00	ng	94
27) 2,4-Dimethylphenol	9.03	122	119674	42.37	ng	95
28) bis(2-Chloroethoxy)methane	9.14	93	166345	39.84	ng	97
29) 2,4-Dichlorophenol	9.30	162	105006	39.29	ng	99
30) 1,2,4-Trichlorobenzene	9.37	180	116082	38.03	ng	98
31) Naphthalene	9.48	128	399008	37.65	ng	100
32) Benzoic acid	9.34	122	32665	18.41	ng	99
33) 4-Chloroaniline	9.63	127	49534	12.45	ng	95
34) Hexachlorobutadiene	9.70	225	62214	35.39	ng	98
35) Caprolactam	10.26	113	33895m	43.51	ng	
36) 4-Chloro-3-methylphenol	10.50	107	116154	37.39	ng	98
37) 2-Methylnaphthalene	10.60	142	257319	39.49	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.88	216	99204	29.19	ng	100
40) Hexachlorocyclopentadiene	10.86	237	55006	50.53	ng	99

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080916\
 Data File : BF089698.D
 Acq On : 9 Aug 2016 20:30
 Operator : UM/SJ
 Sample : PB92711BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 PB92711BS

Manual Integrations
 APPROVED

sohil
 8/10/2016 9:23:10 AM

Quant Time: Aug 10 02:08:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 01:39:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.11	196	68812	41.05	nq	99
43) 2,4,5-Trichlorophenol	11.20	196	67838	37.87	nq	97
45) 1,1'-Biphenyl	11.37	154	274033	33.60	nq	99
46) 2-Chloronaphthalene	11.38	162	236013	38.60	nq	99
47) 2-Nitroaniline	11.60	65	74643	37.88	nq	96
48) Acenaphthylene	12.03	152	378860	37.62	nq	100
49) Dimethylphthalate	11.93	163	288392	40.69	nq	99
50) 2,6-Dinitrotoluene	12.01	165	60456	42.90	nq	# 78
51) Acenaphthene	12.32	154	221175	38.02	nq	99
52) 3-Nitroaniline	12.29	138	44066	26.15	nq	99
53) 2,4-Dinitrophenol	12.51	184	11826	28.03	nq	# 79
54) Dibenzofuran	12.61	168	339141	42.41	nq	99
55) 4-Nitrophenol	12.70	139	48019m	49.34	nq	
56) 2,4-Dinitrotoluene	12.67	165	82210	41.48	nq	# 83
57) Fluorene	13.15	166	271651	39.74	nq	99
58) 2,3,4,6-Tetrachlorophenol	12.85	232	51470	37.84	nq	100
59) Diethylphthalate	13.07	149	290334	39.61	nq	98
60) 4-Chlorophenyl-phenylether	13.19	204	120904	38.57	nq	97
61) 4-Nitroaniline	13.28	138	52203	34.45	nq	91
62) Azobenzene	13.45	77	317723	44.82	nq	98
64) 4,6-Dinitro-2-methylphenol	13.34	198	18403	21.35	nq	98
65) n-Nitrosodiphenylamine	13.41	169	229484	40.67	nq	99
66) 4-Bromophenyl-phenylether	13.98	248	66843	41.37	nq	95
67) Hexachlorobenzene	14.05	284	66437	37.38	nq	# 82
68) Atrazine	14.33	200	71329	45.28	nq	97
69) Pentachlorophenol	14.41	266	33398	45.75	nq	99
70) Phenanthrene	14.70	178	368914	40.61	nq	99
71) Anthracene	14.79	178	391349	40.18	nq	99
72) Carbazole	15.07	167	329403	40.58	nq	99
73) Di-n-butylphthalate	15.67	149	449063	40.91	nq	99
74) Fluoranthene	16.47	202	368097	39.42	nq	98
76) Benzidine	16.71	184	70252	22.76	nq	99
77) Pyrene	16.77	202	339350	37.70	nq	98
79) Butylbenzylphthalate	17.70	149	173216	45.21	nq	98
80) Benzo(a)anthracene	18.35	228	241223	41.23	nq	99
81) 3,3'-Dichlorobenzidine	18.34	252	57419	31.15	nq	97
82) Chrysene	18.40	228	244345	39.81	nq	98
83) Bis(2-ethylhexyl)phthalate	18.45	149	227400	42.87	nq	100
84) Di-n-octyl phthalate	19.22	149	343232	44.97	nq	98
85) Indeno(1,2,3-cd)pyrene	21.20	276	259515	55.68	nq	99
87) Benzo(b)fluoranthene	19.62	252	199756m	34.04	nq	
88) Benzo(k)fluoranthene	19.66	252	235602m	39.79	nq	
89) Benzo(a)pyrene	19.98	252	221520	40.31	nq	98
90) Dibenzo(a,h)anthracene	21.21	278	215296	41.94	nq	97
91) Benzo(g,h,i)perylene	21.54	276	220754	42.01	nq	95

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

