

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079889.D
 Acq On : 11 Jun 2015 12:10
 Operator : TP/IZ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

apatel
 6/13/2015 8:43:00 AM

Quant Time: Jun 12 08:48:18 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 00:59:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	47833	20.00	ng	0.00
21) Naphthalene-d8	8.68	136	186053	20.00	ng	0.00
38) Acenaphthene-d10	10.46	164	94264	20.00	ng	0.00
63) Phenanthrene-d10	11.98	188	199682	20.00	ng	0.00
75) Chrysene-d12	15.07	240	219262	20.00	ng	0.08
86) Perylene-d12	17.04	264	204611	20.00	ng	0.11

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	6.00	112	13028	4.56	ng	0.00
7) Phenol-d6	6.99	99	16005	4.29	ng	0.00
23) Nitrobenzene-d5	7.94	82	11116	3.55	ng	-0.01
41) 2,4,6-Tribromophenol	11.26	330	2915	3.46	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	35714	5.34	ng	0.00
78) Terphenyl-d14	13.73	244	49733	5.15	ng	0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.37	88	3397	2.60	ng	# 80
3) Pyridine	4.22	79	7748m	2.27	ng	
4) n-Nitrosodimethylamine	4.08	42	3428	2.01	ng	# 80
6) Aniline	7.05	93	11914	2.21	ng	96
8) 2-Chlorophenol	7.18	128	7674	2.33	ng	99
9) Benzaldehyde	6.94	77	5445	2.54	ng	97
10) Phenol	7.00	94	9384	2.34	ng	100
11) bis(2-Chloroethyl)ether	7.11	93	7751	2.44	ng	98
12) 1,3-Dichlorobenzene	7.34	146	9083	2.42	ng	99
13) 1,4-Dichlorobenzene	7.42	146	9672	2.30	ng	98
14) 1,2-Dichlorobenzene	7.56	146	8319	2.28	ng	98
15) Benzyl Alcohol	7.52	79	6543	2.19	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.66	45	11695	2.41	ng	96
17) 2-Methylphenol	7.62	107	6667	2.38	ng	96
19) n-Nitroso-di-n-propylamine	7.78	70	6052	2.29	ng	# 97
20) 3+4-Methylphenols	7.77	107	8407	2.31	ng	98
22) Acetophenone	7.79	105	10274	2.24	ng	# 95
24) Nitrobenzene	7.96	77	6696	2.09	ng	98
25) Isophorone	8.20	82	16612	2.46	ng	98
27) 2,4-Dimethylphenol	8.31	122	8163	2.70	ng	94
28) bis(2-Chloroethoxy)methane	8.41	93	9402	2.33	ng	97
29) 2,4-Dichlorophenol	8.52	162	5539	2.17	ng	100
30) 1,2,4-Trichlorobenzene	8.63	180	7937	2.42	ng	99
31) Naphthalene	8.71	128	26887	2.70	ng	99
33) 4-Chloroaniline	8.74	127	9470m	2.36	ng	
34) Hexachlorobutadiene	8.83	225	4386	2.44	ng	95
35) Caprolactam	9.06	113	1766	2.28	ng	# 67
36) 4-Chloro-3-methylphenol	9.21	107	6347	2.26	ng	97
37) 2-Methylnaphthalene	9.40	142	17206	2.44	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.58	216	7754	2.38	ng	96
42) 2,4,6-Trichlorophenol	9.68	196	3979	2.00	ng	92
43) 2,4,5-Trichlorophenol	9.71	196	4919	2.36	ng	94
45) 1,1'-Biphenyl	9.86	154	18487	2.38	ng	98
46) 2-Chloronaphthalene	9.90	162	16687	2.61	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
48) Acenaphthylene	10.32	152	27596	2.52	nq	96
49) Dimethylphthalate	10.15	163	19178	2.58	nq	99
51) Acenaphthene	10.49	154	15537	2.45	nq	99
54) Dibenzofuran	10.66	168	23061	2.58	nq	99
57) Fluorene	11.02	166	19607	2.47	nq	98
59) Diethylphthalate	10.86	149	19266	2.64	nq	93
60) 4-Chlorophenyl-phenylether	10.99	204	9214	2.68	nq	94
62) Azobenzene	11.15	77	16987	2.38	nq	98
65) n-Nitrosodiphenylamine	11.11	169	16499	2.47	nq	95
66) 4-Bromophenyl-phenylether	11.50	248	5808	2.49	nq	# 85
67) Hexachlorobenzene	11.58	284	6050	2.49	nq	# 81
68) Atrazine	11.63	200	4330	2.25	nq	96
70) Phenanthrene	12.01	178	31052	2.58	nq	97
71) Anthracene	12.06	178	28750	2.47	nq	99
72) Carbazole	12.22	167	28219	2.57	nq	99
73) Di-n-butylphthalate	12.55	149	28291	2.48	nq	# 98
74) Fluoranthene	13.31	202	36291	2.86	nq	96
76) Benzidine	13.43	184	15228	2.39	nq	98
77) Pyrene	13.58	202	38061	2.80	nq	97
80) Benzo(a)anthracene	15.05	228	45986	3.53	nq	100
81) 3,3'-Dichlorobenzidine	14.99	252	11076	2.69	nq	# 97
82) Chrysene	15.10	228	48787	3.99	nq	99
83) Bis(2-ethylhexyl)phthalate	15.04	149	14604	2.03	nq	98
84) Di-n-octyl phthalate	15.90	149	22989	2.03	nq	# 84
85) Indeno(1,2,3-cd)pyrene	18.97	276	137042	10.52	nq	98
87) Benzo(b)fluoranthene	16.48	252	55177	4.54	nq	99
88) Benzo(k)fluoranthene	16.51	252	67569	5.34	nq	99
89) Benzo(a)pyrene	16.96	252	71810	6.25	nq	# 96
90) Dibenzo(a,h)anthracene	18.99	278	123670	11.64	nq	98
91) Benzo(g,h,i)perylene	19.54	276	112911	10.05	nq	95

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

