

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052417\
 Data File : BF095397.D
 Acq On : 24 May 2017 19:31
 Operator : SJ/MA
 Sample : I3241-17MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-WC2(0-6)MS

Manual Integrations
 APPROVED

mohammad
 5/25/2017 8:52:42 AM

Quant Time: May 25 04:17:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	83637	20.00	ng	-0.02
21) Naphthalene-d8	9.77	136	331965	20.00	ng	-0.02
38) Acenaphthene-d10	12.59	164	122743	20.00	ng	-0.02
63) Phenanthrene-d10	15.00	188	188473	20.00	ng	-0.02
75) Chrysene-d12	18.69	240	155200	20.00	ng	0.00
86) Perylene-d12	20.37	264	115743	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	545133	108.67	ng	-0.01
7) Phenol-d6	7.30	99	630859	104.81	ng	-0.01
23) Nitrobenzene-d5	8.65	82	356530	67.72	ng	-0.02
41) 2,4,6-Tribromophenol	13.90	330	91855	91.38	ng	-0.02
44) 2-Fluorobiphenyl	11.54	172	647199	75.89	ng	-0.02
78) Terphenyl-d14	17.33	244	491148	69.70	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.24	88	74972	30.47	ng	93
3) Pyridine	2.94	79	149808	26.27	ng	99
4) n-Nitrosodimethylamine	2.87	42	65724	27.65	ng	79
6) Aniline	7.26	93	128431	16.90	ng	97
8) 2-Chlorophenol	7.45	128	195732	34.28	ng	92
9) Benzaldehyde	7.08	77	66325	18.05	ng	# 82
10) Phenol	7.32	94	254981	40.20	ng	93
11) bis(2-Chloroethyl)ether	7.38	93	187954	35.89	ng	98
12) 1,3-Dichlorobenzene	7.65	146	226531	34.97	ng	98
13) 1,4-Dichlorobenzene	7.78	146	229521	34.94	ng	98
14) 1,2-Dichlorobenzene	8.01	146	219982	35.88	ng	97
15) Benzyl Alcohol	8.04	79	153641	39.69	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.24	45	250857	33.85	ng	96
17) 2-Methylphenol	8.26	107	166004	35.53	ng	93
18) Hexachloroethane	8.52	117	56584	25.43	ng	91
19) n-Nitroso-di-n-propylamine	8.45	70	139116	34.17	ng	91
20) 3+4-Methylphenols	8.52	107	206632	32.54	ng	# 56
22) Acetophenone	8.42	105	273265	34.31	ng	# 81
24) Nitrobenzene	8.68	77	191232	34.66	ng	93
25) Isophorone	9.08	82	339058	36.07	ng	96
26) 2-Nitrophenol	9.20	139	78070	26.22	ng	98
27) 2,4-Dimethylphenol	9.33	122	165085	34.29	ng	96
28) bis(2-Chloroethoxy)methane	9.44	93	228051	35.07	ng	97
29) 2,4-Dichlorophenol	9.64	162	147047	32.35	ng	98
30) 1,2,4-Trichlorobenzene	9.70	180	166272	34.14	ng	98
31) Naphthalene	9.80	128	671288	40.23	ng	99
33) 4-Chloroaniline	9.99	127	47529	7.64	ng	# 92
34) Hexachlorobutadiene	10.01	225	81717	31.80	ng	98
35) Caprolactam	10.56	113	39819m	29.17	ng	
36) 4-Chloro-3-methylphenol	10.82	107	148206	30.50	ng	90
37) 2-Methylnaphthalene	10.92	142	408808	37.33	ng	98
39) 1,2,4,5-Tetrachlorobenzene	11.21	216	135200	35.82	ng	99
42) 2,4,6-Trichlorophenol	11.44	196	85754	34.03	ng	98
43) 2,4,5-Trichlorophenol	11.55	196	85598	31.60	ng	# 94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	11.69	154	401200	38.82	nq	98
46) 2-Chloronaphthalene	11.71	162	295449	35.41	nq	95
47) 2-Nitroaniline	11.94	65	84813	37.13	nq	# 79
48) Acenaphthylene	12.37	152	535783	40.99	nq	99
49) Dimethylphthalate	12.24	163	411541	40.84	nq	99
50) 2,6-Dinitrotoluene	12.34	165	58898	28.65	nq	# 92
51) Acenaphthene	12.65	154	344679	44.15	nq	98
52) 3-Nitroaniline	12.62	138	63175	28.63	nq	98
54) Dibenzofuran	12.94	168	448969	41.56	nq	96
55) 4-Nitrophenol	13.07	139	53908	40.68	nq	# 84
56) 2,4-Dinitrotoluene	13.02	165	67798	25.67	nq	96
57) Fluorene	13.49	166	408547	43.49	nq	99
58) 2,3,4,6-Tetrachlorophenol	13.19	232	55572	32.60	nq	95
59) Diethylphthalate	13.38	149	320248	33.16	nq	98
60) 4-Chlorophenyl-phenylether	13.51	204	135300	32.77	nq	92
61) 4-Nitroaniline	13.62	138	57856	28.56	nq	95
62) Azobenzene	13.77	77	263050	33.89	nq	89
65) n-Nitrosodiphenylamine	13.72	169	254261	37.17	nq	98
66) 4-Bromophenyl-phenylether	14.30	248	70411	35.36	nq	99
67) Hexachlorobenzene	14.39	284	67255	32.96	nq	# 86
68) Atrazine	14.64	200	71764	37.16	nq	97
69) Pentachlorophenol	14.77	266	42401	48.45	nq	96
70) Phenanthrene	15.05	178	1802871	166.48	nq	99
71) Anthracene	15.12	178	754577	68.22	nq	97
72) Carbazole	15.41	167	467445	46.44	nq	97
73) Di-n-butylphthalate	15.97	149	458137	36.50	nq	99
74) Fluoranthene	16.81	202	2221966	200.03	nq	99
77) Pvrene	17.11	202	2355774	202.96	nq	98
79) Butylbenzylphthalate	17.99	149	208930	38.70	nq	91
80) Benzo(a)anthracene	18.68	228	1234070	136.63	nq	99
81) 3,3'-Dichlorobenzidine	18.65	252	56922	18.25	nq	98
82) Chrysene	18.73	228	1100905	126.05	nq	98
83) Bis(2-ethylhexyl)phthalate	18.74	149	269120	34.99	nq	# 99
84) Di-n-octyl phthalate	19.50	149	522670	41.44	nq	96
85) Indeno(1,2,3-cd)pvrene	21.70	276	488646	68.89	nq	96
87) Benzo(b)fluoranthene	19.96	252	970890	135.75	nq	# 97
88) Benzo(k)fluoranthene	19.99	252	425008m	61.61	nq	
89) Benzo(a)pyrene	20.31	252	699236	105.95	nq	97
90) Dibenzo(a,h)anthracene	21.70	278	255725	46.92	nq	96
91) Benzo(g,h,i)perylene	22.09	276	459562	85.92	nq	98

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

