

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072816\
 Data File : BF089356.D
 Acq On : 28 Jul 2016 9:51
 Operator : UM/SJ
 Sample : H4155-17MS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 I-4MS

Manual Integrations

sohil
 7/28/2016 5:03:50 PM

Quant Time: Jul 28 13:31:29 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 27 17:12:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	46763	20.00	ng	0.00
21) Naphthalene-d8	7.83	136	200807	20.00	ng	0.00
38) Acenaphthene-d10	9.58	164	96551	20.00	ng	0.00
63) Phenanthrene-d10	11.05	188	184462	20.00	ng	0.00
75) Chrysene-d12	13.67	240	129117	20.00	ng	0.00
86) Perylene-d12	15.01	264	99745	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.14	112	358620	122.45	ng	0.03
7) Phenol-d6	6.22	99	455657	117.73	ng	0.01
23) Nitrobenzene-d5	7.12	82	343081	100.85	ng	0.00
41) 2,4,6-Tribromophenol	10.36	330	110080	137.66	ng	0.00
44) 2-Fluorobiphenyl	8.91	172	658101	100.03	ng	0.00
78) Terphenyl-d14	12.63	244	540009	97.89	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.16	88	45304	33.58	ng	# 83
3) Pyridine	2.74	79	76066m	26.55	ng	
4) n-Nitrosodimethylamine	2.68	42	48051	41.90	ng	95
6) Aniline	6.23	93	73036	17.48	ng	# 1
8) 2-Chlorophenol	6.33	128	154678	46.69	ng	86
9) Benzaldehyde	6.09	77	45748	24.21	ng	95
10) Phenol	6.23	94	177336	41.52	ng	88
11) bis(2-Chloroethyl)ether	6.30	93	187239	51.60	ng	94
12) 1,3-Dichlorobenzene	6.48	146	144006	41.56	ng	96
13) 1,4-Dichlorobenzene	6.56	146	157323	41.14	ng	98
14) 1,2-Dichlorobenzene	6.72	146	150249	41.94	ng	99
15) Benzyl Alcohol	6.71	79	129873	54.07	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.84	45	179104	45.25	ng	90
17) 2-Methylphenol	6.83	107	119692	46.75	ng	96
18) Hexachloroethane	7.05	117	59740	40.75	ng	96
19) n-Nitroso-di-n-propylamine	6.98	70	126695	54.28	ng	93
20) 3+4-Methylphenols	6.99	107	154265	46.91	ng	89
22) Acetophenone	6.97	105	212259	44.40	ng	98
24) Nitrobenzene	7.14	77	175098	45.22	ng	99
25) Isophorone	7.38	82	332905	48.58	ng	98
26) 2-Nitrophenol	7.46	139	79574	45.97	ng	97
27) 2,4-Dimethylphenol	7.51	122	127458	46.22	ng	96
28) bis(2-Chloroethoxy)methane	7.60	93	200062	52.74	ng	99
29) 2,4-Dichlorophenol	7.70	162	127799	50.86	ng	92
30) 1,2,4-Trichlorobenzene	7.78	180	132861	46.34	ng	99
31) Naphthalene	7.85	128	498543m	50.14	ng	
32) Benzoic acid	7.68	122	75184	39.12	ng	97
33) 4-Chloroaniline	7.95	127	39345	10.39	ng	97
34) Hexachlorobutadiene	7.98	225	73033	44.47	ng	96
35) Caprolactam	8.31	113	33932m	44.74	ng	
36) 4-Chloro-3-methylphenol	8.42	107	148573	49.75	ng	94
37) 2-Methylnaphthalene	8.55	142	307906	51.39	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.71	216	115532	34.19	ng	99
40) Hexachlorocyclopentadiene	8.70	237	93303	94.99	ng	97

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42) 2,4,6-Trichlorophenol	8.83	196	83447	51.95	ng	98
43) 2,4,5-Trichlorophenol	8.88	196	105281	51.74	ng #	91
45) 1,1'-Biphenyl	9.00	154	325925	39.65	ng	100
46) 2-Chloronaphthalene	9.03	162	291306	47.18	ng	100
47) 2-Nitroaniline	9.14	65	95831	52.41	ng	94
48) Acenaphthylene	9.44	152	472214	48.56	ng	99
49) Dimethylphthalate	9.32	163	346884	47.21	ng	99
50) 2,6-Dinitrotoluene	9.38	165	73394	48.61	ng #	71
51) Acenaphthene	9.61	154	284897	50.17	ng	99
52) 3-Nitroaniline	9.55	138	43868	27.53	ng	85
53) 2,4-Dinitrophenol	9.67	184	67639	90.35	ng	94
54) Dibenzofuran	9.78	168	403204	52.11	ng	96
55) 4-Nitrophenol	9.75	139	61332m	79.11	ng	
56) 2,4-Dinitrotoluene	9.79	165	108499	54.59	ng #	65
57) Fluorene	10.12	166	344283	50.97	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.91	232	75858	55.84	ng	95
59) Diethylphthalate	10.02	149	390775	52.46	ng	97
60) 4-Chlorophenyl-phenylether	10.12	204	149238	48.17	ng	98
61) 4-Nitroaniline	10.17	138	72363	44.06	ng	97
62) Azobenzene	10.28	77	400323	52.18	ng	85
64) 4,6-Dinitro-2-methylphenol	10.20	198	49788	45.02	ng	90
65) n-Nitrosodiphenylamine	10.25	169	293542	51.00	ng	99
66) 4-Bromophenyl-phenylether	10.60	248	91337	51.86	ng #	89
67) Hexachlorobenzene	10.66	284	85026	47.11	ng	93
68) Atrazine	10.78	200	77457	43.81	ng	97
69) Pentachlorophenol	10.87	266	88805	99.93	ng	98
70) Phenanthrene	11.07	178	478415	50.29	ng	99
71) Anthracene	11.13	178	521562	49.31	ng	100
72) Carbazole	11.29	167	469542	52.73	ng	99
73) Di-n-butylphthalate	11.63	149	583556	48.85	ng #	98
74) Fluoranthene	12.25	202	497072	49.67	ng	97
76) Benzidine	12.41	184	18934	3.97	ng	98
77) Pyrene	12.48	202	497606	50.85	ng	99
79) Butylbenzylphthalate	13.11	149	240193	53.99	ng	95
80) Benzo(a)anthracene	13.66	228	362189	49.71	ng	99
81) 3,3'-Dichlorobenzidine	13.63	252	77168	29.43	ng	96
82) Chrysene	13.70	228	374614	48.30	ng	99
83) Bis(2-ethylhexyl)phthalate	13.67	149	318636	49.54	ng	100
84) Di-n-octyl phthalate	14.27	149	507260	50.55	ng	100
85) Indeno(1,2,3-cd)pyrene	16.21	276	278742	50.55	ng	100
87) Benzo(b)fluoranthene	14.65	252	277367m	44.78	ng	
88) Benzo(k)fluoranthene	14.67	252	328812m	57.74	ng	
89) Benzo(a)pyrene	14.96	252	285524	51.38	ng #	94
90) Dibenzo(a,h)anthracene	16.22	278	235921	53.90	ng	96
91) Benzo(g,h,i)perylene	16.56	276	229403	53.31	ng	97

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

