

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122415\
 Data File : BF084082.D
 Acq On : 24 Dec 2015 16:31
 Operator : UM/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF122415

Manual Integrations
 APPROVED

Sohil
 12/28/2015 1:08:57 PM

Quant Time: Dec 25 03:54:28 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 24 16:48:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	47029	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	189632	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	88516	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	159294	20.00	ng	0.00
75) Chrysene-d12	13.96	240	118317	20.00	ng	0.00
86) Perylene-d12	15.38	264	92999	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	221679	79.92	ng	0.00
7) Phenol-d6	6.45	99	254514	74.15	ng	0.00
23) Nitrobenzene-d5	7.37	82	241101	74.42	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	66402	70.35	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	454822	70.31	ng	0.00
78) Terphenyl-d14	12.90	244	444173	67.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.35	88	45183	39.74	ng	98
3) Pyridine	3.06	79	125064	45.04	ng	98
4) n-Nitrosodimethylamine	3.00	42	49807	40.71	ng	92
6) Aniline	6.46	93	172694	39.80	ng	# 70
8) 2-Chlorophenol	6.59	128	134506	39.05	ng	93
9) Benzaldehyde	6.35	77	68707	37.43	ng	90
10) Phenol	6.46	94	145268	37.00	ng	# 62
11) bis(2-Chloroethyl)ether	6.53	93	106413	37.59	ng	96
12) 1,3-Dichlorobenzene	6.74	146	146571	38.97	ng	96
13) 1,4-Dichlorobenzene	6.82	146	147882	39.08	ng	93
14) 1,2-Dichlorobenzene	6.97	146	131071	37.58	ng	97
15) Benzyl Alcohol	6.96	79	101398	39.01	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.08	45	103068	37.27	ng	58
17) 2-Methylphenol	7.07	107	96196	36.37	ng	# 92
18) Hexachloroethane	7.31	117	52231	37.82	ng	100
19) n-Nitroso-di-n-propylamine	7.22	70	77852	36.82	ng	# 87
20) 3+4-Methylphenols	7.23	107	128949	38.22	ng	# 48
22) Acetophenone	7.22	105	180485	38.09	ng	# 90
24) Nitrobenzene	7.39	77	135789	39.47	ng	92
25) Isophorone	7.63	82	224438	37.45	ng	97
26) 2-Nitrophenol	7.71	139	69294	39.36	ng	93
27) 2,4-Dimethylphenol	7.76	122	119777	40.99	ng	95
28) bis(2-Chloroethoxy)methane	7.84	93	119327	34.15	ng	99
29) 2,4-Dichlorophenol	7.95	162	100379	36.99	ng	95
30) 1,2,4-Trichlorobenzene	8.03	180	116044	39.02	ng	97
31) Naphthalene	8.11	128	388030	38.46	ng	99
32) Benzoic acid	7.88	122	51392	36.68	ng	95
33) 4-Chloroaniline	8.17	127	146242	36.43	ng	# 88
34) Hexachlorobutadiene	8.22	225	62176	36.66	ng	99
35) Caprolactam	8.53	113	29456	39.14	ng	96
36) 4-Chloro-3-methylphenol	8.66	107	120670	39.55	ng	99
37) 2-Methylnaphthalene	8.80	142	227159	35.97	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	107908	38.67	ng	99
40) Hexachlorocyclopentadiene	8.96	237	23338	38.73	ng	98

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42) 2,4,6-Trichlorophenol	9.08	196	68526	37.66	ng	96
43) 2,4,5-Trichlorophenol	9.13	196	70334	38.62	ng #	85
45) 1,1'-Biphenyl	9.26	154	324321	40.56	ng	98
46) 2-Chloronaphthalene	9.29	162	233013	38.33	ng	96
47) 2-Nitroaniline	9.39	65	66181	40.97	ng #	61
48) Acenaphthylene	9.70	152	367606	36.82	ng	99
49) Dimethylphthalate	9.56	163	272832	37.22	ng #	90
50) 2,6-Dinitrotoluene	9.63	165	62213	39.98	ng #	83
51) Acenaphthene	9.87	154	214153	36.88	ng	99
52) 3-Nitroaniline	9.80	138	70265	42.08	ng #	77
53) 2,4-Dinitrophenol	9.92	184	15921	36.77	ng	97
54) Dibenzofuran	10.04	168	283949	35.56	ng #	90
55) 4-Nitrophenol	9.98	139	33377	39.16	ng	96
56) 2,4-Dinitrotoluene	10.03	165	74594	36.31	ng #	80
57) Fluorene	10.38	166	250668	36.93	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	51407	39.60	ng	97
59) Diethylphthalate	10.26	149	251856	33.82	ng	96
60) 4-Chlorophenyl-phenylether	10.37	204	104654	33.43	ng #	78
61) 4-Nitroaniline	10.41	138	62322	40.38	ng	91
62) Azobenzene	10.53	77	228946	37.32	ng	87
64) 4,6-Dinitro-2-methylphenol	10.45	198	33234	37.59	ng	78
65) n-Nitrosodiphenylamine	10.50	169	227694	40.10	ng	98
66) 4-Bromophenyl-phenylether	10.86	248	72667	39.38	ng #	80
67) Hexachlorobenzene	10.93	284	72031	35.65	ng #	68
68) Atrazine	11.02	200	59820	34.98	ng	99
69) Pentachlorophenol	11.14	266	21177	36.52	ng	98
70) Phenanthrene	11.34	178	359963	38.12	ng	100
71) Anthracene	11.40	178	364301	36.89	ng	98
72) Carbazole	11.55	167	303869	36.17	ng	99
73) Di-n-butylphthalate	11.88	149	398002	36.78	ng #	98
74) Fluoranthene	12.53	202	391024	38.36	ng	95
76) Benzidine	12.65	184	147279	36.62	ng	98
77) Pyrene	12.76	202	386583	37.87	ng	100
79) Butylbenzylphthalate	13.38	149	161318	38.18	ng #	80
80) Benzo(a)anthracene	13.95	228	274476	37.71	ng	99
81) 3,3'-Dichlorobenzidine	13.91	252	87179	39.55	ng #	96
82) Chrysene	13.99	228	269150m	40.10	ng	
83) Bis(2-ethylhexyl)phthalate	13.94	149	208595	38.21	ng #	97
84) Di-n-octyl phthalate	14.55	149	298415	38.84	ng	100
85) Indeno(1,2,3-cd)pyrene	16.77	276	251541	37.43	ng	99
87) Benzo(b)fluoranthene	14.98	252	258073m	37.37	ng	
88) Benzo(k)fluoranthene	15.00	252	189371m	40.46	ng	
89) Benzo(a)pyrene	15.32	252	208212	36.59	ng	99
90) Dibenzo(a,h)anthracene	16.77	278	209481	37.73	ng #	95
91) Benzo(g,h,i)perylene	17.19	276	212150	37.43	ng	98

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

