

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012916\
 Data File : BF084662.D
 Acq On : 30 Jan 2016 3:46
 Operator : SJ/IZ
 Sample : H1261-11MS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SCARBOROUGH1516-00(0-4)MS

Manual Integrations
 APPROVED

mohammad
 2/1/2016 2:24:33 PM

Quant Time: Feb 01 07:36:17 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:54:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	170956	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	668380	20.00	ng	-0.01
38) Acenaphthene-d10	9.77	164	317423	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	590820	20.00	ng	0.00
75) Chrysene-d12	13.87	240	409650	20.00	ng	0.00
86) Perylene-d12	15.25	264	276099	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	1344960	118.76	ng	0.02
7) Phenol-d6	6.38	99	1645672	121.24	ng	0.01
23) Nitrobenzene-d5	7.30	82	1164243	97.35	ng	0.00
41) 2,4,6-Tribromophenol	10.56	330	484832	145.73	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1864632	87.31	ng	0.00
78) Terphenyl-d14	12.83	244	1856693	102.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	187387	35.48	ng	# 78
3) Pyridine	3.06	79	259542m	18.45	ng	
4) n-Nitrosodimethylamine	2.98	42	264544	38.46	ng	# 84
6) Aniline	6.40	93	67868	3.90	ng	# 1
8) 2-Chlorophenol	6.51	128	508738	40.50	ng	83
10) Phenol	6.40	94	596751	39.45	ng	82
11) bis(2-Chloroethyl)ether	6.48	93	542489	46.11	ng	96
12) 1,3-Dichlorobenzene	6.67	146	521756	38.77	ng	96
13) 1,4-Dichlorobenzene	6.74	146	514862	38.98	ng	98
14) 1,2-Dichlorobenzene	6.90	146	530742	43.64	ng	98
15) Benzyl Alcohol	6.88	79	367094	39.73	ng	91
16) 2,2'-oxybis(1-Chloropropan	7.01	45	798788	39.67	ng	91
17) 2-Methylphenol	7.00	107	429360	44.09	ng	93
18) Hexachloroethane	7.24	117	212446	41.11	ng	# 86
19) n-Nitroso-di-n-propylamine	7.15	70	361953	43.65	ng	# 74
20) 3+4-Methylphenols	7.15	107	518510	43.88	ng	# 66
22) Acetophenone	7.14	105	748300	42.87	ng	# 99
24) Nitrobenzene	7.31	77	509265	40.06	ng	94
25) Isophorone	7.56	82	1043319	46.56	ng	95
26) 2-Nitrophenol	7.63	139	274403	45.39	ng	# 84
27) 2,4-Dimethylphenol	7.68	122	453634	42.82	ng	95
28) bis(2-Chloroethoxy)methane	7.77	93	612637	45.56	ng	99
29) 2,4-Dichlorophenol	7.88	162	451525	48.73	ng	95
30) 1,2,4-Trichlorobenzene	7.95	180	455753	43.02	ng	96
31) Naphthalene	8.03	128	1649568	48.89	ng	100
32) Benzoic acid	7.81	122	319944	39.35	ng	95
34) Hexachlorobutadiene	8.16	225	274594	43.86	ng	99
35) Caprolactam	8.46	113	126293m	47.50	ng	
36) 4-Chloro-3-methylphenol	8.59	107	496440	48.67	ng	93
37) 2-Methylnaphthalene	8.73	142	1109260	53.56	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.89	216	401724	39.02	ng	99
40) Hexachlorocyclopentadiene	8.88	237	433519	92.62	ng	99
42) 2,4,6-Trichlorophenol	9.01	196	316891	49.55	ng	95
43) 2,4,5-Trichlorophenol	9.06	196	315495	47.79	ng	# 94

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012916\
 Data File : BF084662.D
 Acq On : 30 Jan 2016 3:46
 Operator : SJ/IZ
 Sample : H1261-11MS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SCARBOROUGH1516-00(0-4)MS

Manual Integrations
 APPROVED

mohammad
 2/1/2016 2:24:33 PM

Quant Time: Feb 01 07:36:17 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:54:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	9.20	154	1241069	42.66	nq	95
46) 2-Chloronaphthalene	9.22	162	945369	44.54	nq #	86
47) 2-Nitroaniline	9.31	65	254960	40.00	nq	97
48) Acenaphthylene	9.63	152	1497692	46.57	nq	98
49) Dimethylphthalate	9.50	163	1196809	50.24	nq #	91
50) 2,6-Dinitrotoluene	9.56	165	268404	51.20	nq #	90
51) Acenaphthene	9.80	154	1015677	47.16	nq	97
52) 3-Nitroaniline	9.72	138	22684	4.14	nq	91
53) 2,4-Dinitrophenol	9.84	184	298611	115.45	nq	90
54) Dibenzofuran	9.97	168	1361376	52.97	nq	97
55) 4-Nitrophenol	9.90	139	351977	87.49	nq #	82
56) 2,4-Dinitrotoluene	9.96	165	339609	49.87	nq #	91
57) Fluorene	10.32	166	1049606	48.18	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.09	232	265624	53.21	nq	88
59) Diethylphthalate	10.20	149	1168837	50.12	nq	97
60) 4-Chlorophenyl-phenylether	10.30	204	436190	44.60	nq	93
61) 4-Nitroaniline	10.34	138	170862	32.96	nq	84
62) Azobenzene	10.46	77	1126259	50.46	nq	91
64) 4,6-Dinitro-2-methylphenol	10.36	198	173670	45.00	nq #	53
65) n-Nitrosodiphenylamine	10.43	169	695177	36.29	nq	98
66) 4-Bromophenyl-phenylether	10.80	248	329647	49.54	nq #	94
67) Hexachlorobenzene	10.85	284	320201	44.24	nq #	84
68) Atrazine	10.96	200	137159	23.90	nq	96
69) Pentachlorophenol	11.06	266	356426	101.93	nq	100
70) Phenanthrene	11.26	178	1437537	44.00	nq	97
71) Anthracene	11.32	178	1635699	49.24	nq	98
72) Carbazole	11.47	167	816116	28.51	nq	100
73) Di-n-butylphthalate	11.81	149	1602392	46.02	nq #	96
74) Fluoranthene	12.45	202	1655745	51.26	nq	95
76) Benzidine	12.59	184	34603	2.22	nq #	93
77) Pyrene	12.68	202	1643305	53.19	nq	99
79) Butylbenzylphthalate	13.31	149	744026	56.37	nq #	90
80) Benzo(a)anthracene	13.86	228	1195866	46.83	nq	98
82) Chrysene	13.91	228	1208111	48.00	nq	99
83) Bis(2-ethylhexyl)phthalate	13.87	149	926037	54.08	nq #	97
84) Di-n-octyl phthalate	14.48	149	1463411	54.09	nq #	89
85) Indeno(1,2,3-cd)pyrene	16.57	276	911211	40.95	nq	99
87) Benzo(b)fluoranthene	14.88	252	1196595m	66.93	nq	
88) Benzo(k)fluoranthene	14.90	252	803687m	54.42	nq	
89) Benzo(a)pyrene	15.20	252	863679m	55.85	nq	
90) Dibenzo(a,h)anthracene	16.58	278	775008	55.61	nq	96
91) Benzo(g,h,i)perylene	16.96	276	768801	54.64	nq	98

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

