

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020718\
 Data File : BN000038.D
 Acq On : 07 Feb 2018 15:10
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:37:31 PM

Quant Time: Feb 07 18:23:28 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 14:04:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	485623	20.00	ng	0.00
21) Naphthalene-d8	11.31	136	2410301	20.00	ng	0.00
38) Acenaphthene-d10	15.09	164	1387169	20.00	ng	0.00
63) Phenanthrene-d10	17.81	188	2901265	20.00	ng	0.01
75) Chrysene-d12	21.94	240	2904965	20.00	ng	0.02
86) Perylene-d12	24.43	264	3206337	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.94	112	5055798	153.50	ng	0.00
7) Phenol-d6	7.59	99	7443344	148.78	ng	0.00
23) Nitrobenzene-d5	9.65	82	7001837	175.66	ng	0.00
41) 2,4,6-Tribromophenol	16.56	330	2370311	169.98	ng	0.00
44) 2-Fluorobiphenyl	13.72	172	13936869	134.56	ng	0.00
78) Terphenyl-d14	20.37	244	14188160	128.91	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.66	88	950919	72.981	ng	94
3) Pyridine	4.09	79	3249696	82.237	ng	98
4) n-Nitrosodimethylamine	4.00	42	882939	75.944	ng	# 86
6) Aniline	7.77	93	5122696	75.113	ng	91
8) 2-Chlorophenol	8.02	128	2883840	79.148	ng	86
9) Benzaldehyde	7.58	77	1955410	76.548	ng	87
10) Phenol	7.62	94	3926687	76.787	ng	85
11) bis(2-Chloroethyl)ether	7.87	93	3179493	77.937	ng	# 82
12) 1,3-Dichlorobenzene	8.36	146	3158238	78.077	ng	99
13) 1,4-Dichlorobenzene	8.50	146	3229518	77.738	ng	97
14) 1,2-Dichlorobenzene	8.83	146	3162193	77.040	ng	97
15) Benzyl Alcohol	8.70	79	2823825	76.861	ng	89
16) 2,2'-oxybis(1-Chloropropan	9.00	45	3318149	75.199	ng	79
17) 2-Methylphenol	8.90	107	2873860	77.937	ng	93
18) Hexachloroethane	9.57	117	1187787	83.375	ng	# 81
19) n-Nitroso-di-n-propylamine	9.28	70	2545632	75.509	ng	# 81
20) 3+4-Methylphenols	9.23	107	4054013	78.496	ng	92
22) Acetophenone	9.30	105	5396812	75.809	ng	# 86
24) Nitrobenzene	9.69	77	3750848	87.733	ng	# 91
25) Isophorone	10.22	82	7142749	75.650	ng	95
26) 2-Nitrophenol	10.41	139	1570713	119.956	ng	# 89
27) 2,4-Dimethylphenol	10.45	122	3038383	78.484	ng	94
28) bis(2-Chloroethoxy)methane	10.70	93	4375766	80.348	ng	96
29) 2,4-Dichlorophenol	10.95	162	2847166	85.713	ng	96
30) 1,2,4-Trichlorobenzene	11.17	180	3115337	80.091	ng	97
31) Naphthalene	11.36	128	10139987	75.012	ng	98
32) Benzoic acid	10.61	122	2472065	123.313	ng	87
33) 4-Chloroaniline	11.46	127	4768678	76.091	ng	94
34) Hexachlorobutadiene	11.64	225	1621957	81.847	ng	96
35) Caprolactam	12.25	113	1197810	81.124	ng	91
36) 4-Chloro-3-methylphenol	12.55	107	3548714	78.570	ng	97
37) 2-Methylnaphthalene	12.94	142	7162712	74.985	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	3322509	82.230	ng	# 96
40) Hexachlorocyclopentadiene	13.27	237	1558610	115.492	ng	93

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020718\
 Data File : BN000038.D
 Acq On : 07 Feb 2018 15:10
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:37:31 PM

Quant Time: Feb 07 18:23:28 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 14:04:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.52	196	2154566	90.216	nq	97
43) 2,4,5-Trichlorophenol	13.59	196	2511565	89.301	nq #	94
45) 1,1'-Biphenyl	13.93	154	9330181	75.279	nq	94
46) 2-Chloronaphthalene	13.97	162	7231560	77.719	nq	95
47) 2-Nitroaniline	14.16	65	2223370	101.175	nq #	79
48) Acenaphthylene	14.81	152	11651360	74.571	nq	97
49) Dimethylphthalate	14.53	163	9137302	73.702	nq	98
50) 2,6-Dinitrotoluene	14.65	165	2043557	92.377	nq #	89
51) Acenaphthene	15.15	154	7525348	76.327	nq	98
52) 3-Nitroaniline	14.98	138	2452339	90.430	nq #	87
53) 2,4-Dinitrophenol	15.17	184	701091	137.830	nq #	1
54) Dibenzofuran	15.47	168	10225372	72.271	nq	92
55) 4-Nitrophenol	15.25	139	1888116	93.847	nq #	83
56) 2,4-Dinitrotoluene	15.42	165	2811493	97.433	nq #	90
57) Fluorene	16.12	166	8235510	69.129	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.69	232	1982622m	86.562	nq	
59) Diethylphthalate	15.87	149	9372236	72.353	nq	99
60) 4-Chlorophenyl-phenylether	16.10	204	3885521	73.794	nq	92
61) 4-Nitroaniline	16.13	138	2468327	85.398	nq	90
62) Azobenzene	16.40	77	9228297	71.291	nq	91
64) 4,6-Dinitro-2-methylphenol	16.18	198	1225791	136.867	nq	87
65) n-Nitrosodiphenylamine	16.32	169	7532369	79.449	nq	99
66) 4-Bromophenyl-phenylether	17.00	248	2283897	84.125	nq #	91
67) Hexachlorobenzene	17.12	284	2617546	83.688	nq #	92
68) Atrazine	17.25	200	2538813	87.848	nq	95
69) Pentachlorophenol	17.44	266	1771299	94.557	nq	98
70) Phenanthrene	17.85	178	12570448	72.008	nq	94
71) Anthracene	17.94	178	12585341	72.775	nq	95
72) Carbazole	18.20	167	12624254	70.557	nq #	95
73) Di-n-butylphthalate	18.73	149	13088854	65.873	nq #	89
74) Fluoranthene	19.83	202	13156104	66.373	nq	81
76) Benzidine	19.99	184	8157422	110.936	nq #	93
77) Pyrene	20.19	202	13771073	75.141	nq #	86
79) Butylbenzylphthalate	21.04	149	7248467	99.222	nq	97
80) Benzo(a)anthracene	21.92	228	13475051	74.817	nq #	91
81) 3,3'-Dichlorobenzidine	21.84	252	5376328	89.078	nq #	96
82) Chrysene	21.98	228	12926549	72.266	nq #	90
83) Bis(2-ethylhexyl)phthalate	21.82	149	10322735	85.396	nq #	99
84) Di-n-octyl phthalate	22.78	149	15112473	77.243	nq	95
85) Indeno(1,2,3-cd)pyrene	22.03	276	18159879	78.704	nq #	86
87) Benzo(b)fluoranthene	23.68	252	14702362	77.119	nq #	93
88) Benzo(k)fluoranthene	23.73	252	14289130	74.755	nq #	90
89) Benzo(a)pyrene	24.33	252	14394820	77.725	nq #	93
90) Dibenzo(a,h)anthracene	27.04	278	14905136	76.891	nq #	87
91) Benzo(g,h,i)perylene	27.83	276	15451758	78.431	nq #	86

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

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020718\
Data File : BN000038.D
Acq On : 07 Feb 2018 15:10
Operator : JU/SJ
Sample : SSTDICC080
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
Client Sampled :
SSTDICC080

Manual Integrations
APPROVED

Sohil
2/8/2018 3:37:31 PM

Quant Time: Feb 07 18:23:28 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 07 14:04:10 2018
Response via : Initial Calibration

