

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020618\
 Data File : BN000026.D
 Acq On : 06 Feb 2018 11:19
 Operator :
 Sample : MDL-S-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-S-05

Manual Integrations
 APPROVED

Sohil
 2/7/2018 5:15:31 PM

Quant Time: Feb 07 16:42:13 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 15:24:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.48	152	333108	20.00	ng	-0.01
21) Naphthalene-d8	11.32	136	1628274	20.00	ng	-0.01
38) Acenaphthene-d10	15.08	164	984350	20.00	ng	-0.01
63) Phenanthrene-d10	17.80	188	2342383	20.00	ng	-0.01
75) Chrysene-d12	21.94	240	2679407	20.00	ng	-0.01
86) Perylene-d12	24.44	264	2792716	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.95	112	3302746	146.19	ng	0.00
7) Phenol-d6	7.59	99	4642983	135.30	ng	0.00
23) Nitrobenzene-d5	9.65	82	2293216	85.16	ng	-0.01
41) 2,4,6-Tribromophenol	16.55	330	1255437	126.87	ng	0.00
44) 2-Fluorobiphenyl	13.72	172	6158877	83.80	ng	-0.01
78) Terphenyl-d14	20.37	244	9306804	91.68	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.66	88	27362	3.061	ng	# 85
3) Pyridine	4.14	79	47547	1.754	ng	# 89
4) n-Nitrosodimethylamine	4.02	42	21143	2.651	ng	# 96
6) Aniline	7.78	93	116398	2.488	ng	# 90
8) 2-Chlorophenol	8.02	128	86778	3.472	ng	# 91
9) Benzaldehyde	7.59	77	73654	1.831	ng	# 98
10) Phenol	7.62	94	126689	3.612	ng	# 93
11) bis(2-Chloroethyl)ether	7.88	93	82897	2.962	ng	# 82
12) 1,3-Dichlorobenzene	8.36	146	81869	2.951	ng	# 96
13) 1,4-Dichlorobenzene	8.51	146	85354	2.995	ng	# 98
14) 1,2-Dichlorobenzene	8.83	146	83367	2.961	ng	# 90
15) Benzyl Alcohol	8.70	79	60622	2.406	ng	# 90
16) 2,2'-oxybis(1-Chloropropan	9.00	45	90386	2.986	ng	# 78
17) 2-Methylphenol	8.90	107	71005	2.807	ng	# 97
18) Hexachloroethane	9.58	117	26554	2.717	ng	# 81
19) n-Nitroso-di-n-propylamine	9.28	70	48768	2.109	ng	# 85
20) 3+4-Methylphenols	9.23	107	91917	2.595	ng	# 91
22) Acetophenone	9.30	105	130411	2.712	ng	# 85
24) Nitrobenzene	9.69	77	101273	3.506	ng	# 94
25) Isophorone	10.22	82	145091	2.275	ng	# 94
26) 2-Nitrophenol	10.41	139	18796	8.987	ng	# 86
27) 2,4-Dimethylphenol	10.45	122	97310	3.721	ng	# 93
28) bis(2-Chloroethoxy)methane	10.70	93	112621	3.061	ng	# 95
29) 2,4-Dichlorophenol	10.95	162	57450	2.560	ng	# 98
30) 1,2,4-Trichlorobenzene	11.17	180	77112	2.935	ng	# 97
31) Naphthalene	11.36	128	281685	3.085	ng	# 98
32) Benzoic acid	10.48	122	20578m	7.789	ng	# 96
33) 4-Chloroaniline	11.46	127	105917	2.502	ng	# 97
34) Hexachlorobutadiene	11.65	225	37962	2.836	ng	# 94
35) Caprolactam	12.18	113	15360	1.540	ng	# 94
36) 4-Chloro-3-methylphenol	12.54	107	69763	2.286	ng	# 94
37) 2-Methylnaphthalene	12.94	142	192426	2.982	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	86918	3.031	ng	# 95
40) Hexachlorocyclopentadiene	13.28	237	27148	2.835	ng	# 93

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020618\
 Data File : BN000026.D
 Acq On : 06 Feb 2018 11:19
 Operator :
 Sample : MDL-S-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampled :
 MDL-S-05

Manual Integrations
 APPROVED

Sohil
 2/7/2018 5:15:31 PM

Quant Time: Feb 07 16:42:13 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 15:24:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.52	196	34435	2.032	nq	98
43) 2,4,5-Trichlorophenol	13.59	196	43087	2.159	nq	# 91
45) 1,1'-Biphenyl	13.93	154	287111	3.264	nq	98
46) 2-Chloronaphthalene	13.97	162	198352	3.004	nq	97
47) 2-Nitroaniline	14.16	65	29204	8.647	nq	96
48) Acenaphthylene	14.81	152	276602	2.495	nq	98
49) Dimethylphthalate	14.52	163	228995	2.603	nq	99
50) 2,6-Dinitrotoluene	14.64	165	27712	7.007	nq	92
51) Acenaphthene	15.15	154	205221	2.933	nq	98
52) 3-Nitroaniline	14.97	138	33396	7.453	nq	98
53) 2,4-Dinitrophenol	15.17	184	5935	6.136	nq	# 39
54) Dibenzofuran	15.47	168	326324	3.250	nq	95
55) 4-Nitrophenol	15.24	139	22121	8.156	nq	# 85
56) 2,4-Dinitrotoluene	15.42	165	32090	8.319	nq	# 93
57) Fluorene	16.12	166	251475	2.975	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.69	232	35797	2.202	nq	# 83
59) Diethylphthalate	15.86	149	224566	2.443	nq	96
60) 4-Chlorophenyl-phenylether	16.10	204	117371	3.141	nq	90
61) 4-Nitroaniline	16.11	138	35504	7.475	nq	97
62) Azobenzene	16.39	77	228567	2.488	nq	92
64) 4,6-Dinitro-2-methylphenol	16.17	198	9982	9.905	nq	# 82
65) n-Nitrosodiphenylamine	16.31	169	207045	2.705	nq	98
66) 4-Bromophenyl-phenylether	16.99	248	62035	2.830	nq	# 82
67) Hexachlorobenzene	17.10	284	74917	2.967	nq	# 82
68) Atrazine	17.23	200	52669	2.257	nq	95
69) Pentachlorophenol	17.44	266	22138	1.464	nq	92
70) Phenanthrene	17.85	178	448582	3.183	nq	98
71) Anthracene	17.93	178	391917	2.807	nq	98
72) Carbazole	18.19	167	378090	2.617	nq	98
73) Di-n-butylphthalate	18.73	149	281438	1.754	nq	# 96
74) Fluoranthene	19.82	202	434564	2.715	nq	95
76) Benzidine	19.99	184	88557	1.306	nq	95
77) Pyrene	20.18	202	478070	2.828	nq	97
79) Butylbenzylphthalate	21.05	149	89485	6.135	nq	# 89
80) Benzo(a)anthracene	21.92	228	500245	3.011	nq	97
81) 3,3'-Dichlorobenzidine	21.84	252	110270	1.981	nq	# 95
82) Chrysene	21.97	228	523152	3.171	nq	97
83) Bis(2-ethylhexyl)phthalate	21.83	149	140830	1.263	nq	# 93
84) Di-n-octyl phthalate	22.80	149	209979	1.164	nq	# 99
85) Indeno(1,2,3-cd)pyrene	27.02	276	533967	2.509	nq	# 86
87) Benzo(b)fluoranthene	23.68	252	464135	2.795	nq	# 94
88) Benzo(k)fluoranthene	23.73	252	491388	2.951	nq	# 93
89) Benzo(a)pyrene	24.33	252	423629	2.626	nq	# 95
90) Dibenzo(a,h)anthracene	27.03	278	453665	2.687	nq	# 90
91) Benzo(g,h,i)perylene	27.80	276	457780	2.668	nq	# 86

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

```
Data Path : Z:\HPCHEM1\BNA N\DATA\BN020618\
Data File : BN000026.D
Acq On    : 06 Feb 2018  11:19
Operator  :
Sample    : MDL-S-05
Misc      :
ALS Vial  : 6      Sample Multiplier: 1
```

Instrument :
BNA_N
ClientSampleId :
MDL-S-05

Manual Integrations APPROVED

Sohil
2/7/2018 5:15:31 PM

Quant Time: Feb 07 16:42:13 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 06 15:24:10 2018
Response via : Initial Calibration

