

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM043019\
 Data File : BM020082.D
 Acq On : 30 Apr 2019 19:11
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 5/2/2019 2:20:59 AM

Quant Time: May 01 03:54:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:29:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	95149	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	376539	20.00	ng	0.00
39) Acenaphthene-d10	14.44	164	231311	20.00	ng	0.00
64) Phenanthrene-d10	17.18	188	552215	20.00	ng	0.00
76) Chrysene-d12	21.37	240	586159	20.00	ng	0.00
87) Perylene-d12	23.65	264	643204	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	962761	163.98	ng	0.00
7) Phenol-d6	6.96	99	1308447	148.08	ng	0.00
23) Nitrobenzene-d5	8.96	82	1564336	185.79	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	584026	156.85	ng	0.00
45) 2-Fluorobiphenyl	13.06	172	3003004	161.79	ng	0.00
79) Terphenyl-d14	19.81	244	5459043	164.27	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.31	88	231545	91.391 ng	97
3) Pyridine	3.70	79	575964	81.175 ng	99
4) n-Nitrosodimethylamine	3.62	42	208278	91.123 ng	88
6) Aniline	7.13	93	821697	72.001 ng	100
8) 2-Chlorophenol	7.36	128	516705	70.761 ng	98
9) Benzaldehyde	6.95	77	393371	76.293 ng	99
10) Phenol	6.98	94	703345	72.857 ng	98
11) bis(2-Chloroethyl)ether	7.23	93	513131	72.648 ng	100
12) 1,3-Dichlorobenzene	7.68	146	595253	77.649 ng	97
13) 1,4-Dichlorobenzene	7.83	146	599330	77.056 ng	95
14) 1,2-Dichlorobenzene	8.14	146	574105	74.443 ng	98
15) Benzyl Alcohol	8.03	79	626423	77.984 ng	100
16) 2,2'-oxybis(1-Chloropropan	8.33	45	399686	59.360 ng	97
17) 2-Methylphenol	8.23	107	440598	68.959 ng	98
18) Hexachloroethane	8.86	117	254517	84.184 ng	95
19) n-Nitroso-di-n-propylamine	8.60	70	539832	69.503 ng	97
20) 3+4-Methylphenols	8.56	107	602849	68.685 ng	99
22) Acetophenone	8.62	105	826505	80.589 ng	97
24) Nitrobenzene	9.00	77	796266	91.292 ng	98
25) Isophorone	9.52	82	1321787	80.646 ng	99
26) 2-Nitrophenol	9.70	139	261115	70.250 ng	98
27) 2,4-Dimethylphenol	9.76	122	400955	76.293 ng	97
28) bis(2-Chloroethoxy)methane	10.00	93	727472	79.264 ng	99
29) 2,4-Dichlorophenol	10.23	162	515309	86.334 ng	99
30) 1,2,4-Trichlorobenzene	10.45	180	620553	96.597 ng	100
31) Naphthalene	10.64	128	1558832	77.183 ng	99
32) Benzoic acid	9.91	122	295119m	63.591 ng	
33) 4-Chloroaniline	10.75	127	644950	77.517 ng	98
34) Hexachlorobutadiene	10.90	225	436131	115.109 ng	99
35) Caprolactam	11.56	113	155547m	68.847 ng	
36) 4-Chloro-3-methylphenol	11.87	107	586929	79.139 ng	97
37) 2-Methylnaphthalene	12.25	142	1138628	77.438 ng	98
38) 1-Methylnaphthalene	12.47	142	1106665	76.227 ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.62	216	727655	101.972 ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM043019\
 Data File : BM020082.D
 Acq On : 30 Apr 2019 19:11
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 5/2/2019 2:20:59 AM

Quant Time: May 01 03:54:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:29:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.59	237	441900	230.267	ng	98
43) 2,4,6-Trichlorophenol	12.86	196	428197	88.510	ng	100
44) 2,4,5-Trichlorophenol	12.93	196	466739	92.695	ng	98
46) 1,1'-Biphenyl	13.27	154	1513456	75.543	ng	98
47) 2-Chloronaphthalene	13.31	162	1201667	78.634	ng	98
48) 2-Nitroaniline	13.52	65	452555	82.158	ng	99
49) Acenaphthylene	14.16	152	1938554	72.908	ng	99
50) Dimethylphthalate	13.90	163	1532832	75.128	ng	100
51) 2,6-Dinitrotoluene	14.02	165	334317	74.100	ng	99
52) Acenaphthene	14.50	154	1103928	74.880	ng	100
53) 3-Nitroaniline	14.35	138	318242	70.642	ng	94
54) 2,4-Dinitrophenol	14.56	184	161729	70.867	ng	95
55) Dibenzofuran	14.84	168	1855617	76.971	ng	98
56) 4-Nitrophenol	14.65	139	272616	87.298	ng	97
57) 2,4-Dinitrotoluene	14.81	165	471023	72.629	ng	96
58) Fluorene	15.49	166	1540959	75.434	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.06	232	443066	96.438	ng	99
60) Diethylphthalate	15.26	149	1543666	72.746	ng	99
61) 4-Chlorophenyl-phenylether	15.48	204	871915	88.651	ng	97
62) 4-Nitroaniline	15.52	138	353081	79.548	ng	99
63) Azobenzene	15.77	77	1827938	80.597	ng	99
65) 4,6-Dinitro-2-methylphenol	15.56	198	235393	66.810	ng	96
66) n-Nitrosodiphenylamine	15.70	169	1303456	72.200	ng	98
67) 4-Bromophenyl-phenylether	16.37	248	547078	86.169	ng	97
68) Hexachlorobenzene	16.47	284	617854	81.105	ng	97
69) Atrazine	16.65	200	498012	81.260	ng	98
70) Pentachlorophenol	16.82	266	382674	98.314	ng	98
71) Phenanthrene	17.23	178	2432626	74.233	ng	99
72) Anthracene	17.32	178	2448382	75.057	ng	99
73) Carbazole	17.59	167	2203629	72.355	ng	99
74) Di-n-butylphthalate	18.15	149	2689832	67.251	ng	98
75) Fluoranthene	19.24	202	3076845	81.587	ng	99
77) Benzidine	19.43	184	1833021	113.339	ng	99
78) Pyrene	19.60	202	3162949	81.483	ng	100
80) Butylbenzylphthalate	20.50	149	1219818	64.818	ng	94
81) Benzo(a)anthracene	21.35	228	3318431	89.370	ng	99
82) 3,3'-Dichlorobenzidine	21.29	252	1286850	93.149	ng	99
83) Chrysene	21.40	228	3188342	89.536	ng	99
84) Bis(2-ethylhexyl)phthalate	21.27	149	1839898	65.670	ng	100
85) Di-n-octyl phthalate	22.17	149	3123738	68.713	ng	100
86) Indeno(1,2,3-cd)pyrene	25.99	276	3897629	111.266	ng	# 100
88) Benzo(b)fluoranthene	22.96	252	3385795	79.275	ng	100
89) Benzo(k)fluoranthene	23.01	252	3173549	78.927	ng	99
90) Benzo(a)pyrene	23.56	252	3139910	82.972	ng	99
91) Dibenzo(a,h)anthracene	26.00	278	3293016	90.255	ng	99
92) Benzo(g,h,i)perylene	26.71	276	3109146	94.084	ng	99

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

