

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016487.D
 Acq On : 21 Aug 2018 15:46
 Operator : SJ/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:06:57 PM

Quant Time: Aug 21 16:27:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 14:43:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	152961	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	627210	20.00	ng	0.00
38) Acenaphthene-d10	14.67	164	429106	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	1262108	20.00	ng	0.00
75) Chrysene-d12	21.55	240	1953145	20.00	ng	0.00
86) Perylene-d12	23.84	264	1932348	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	1326230m	145.71	ng	0.00
7) Phenol-d6	7.20	99	1742885	142.46	ng	0.00
23) Nitrobenzene-d5	9.22	82	2278893	151.32	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	2199035	339.30	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	8478346	205.45	ng	0.00
78) Terphenyl-d14	19.99	244	12550342	97.12	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.38	88	335447m	73.595	ng	
3) Pyridine	3.80	79	739607m	80.599	ng	
4) n-Nitrosodimethylamine	3.73	42	320474m	30.310	ng	
6) Aniline	7.36	93	945657	74.700	ng	91
8) 2-Chlorophenol	7.59	128	781484	69.643	ng	96
9) Benzaldehyde	7.18	77	469767	58.400	ng	# 79
10) Phenol	7.23	94	847292	71.499	ng	97
11) bis(2-Chloroethyl)ether	7.46	93	631514	70.194	ng	99
12) 1,3-Dichlorobenzene	7.90	146	995381	77.396	ng	# 85
13) 1,4-Dichlorobenzene	8.06	146	1030182	79.238	ng	98
14) 1,2-Dichlorobenzene	8.38	146	1033038	80.884	ng	98
15) Benzyl Alcohol	8.29	79	772621	69.579	ng	# 82
16) 2,2'-oxybis(1-Chloropropan	8.55	45	406160	35.864	ng	# 54
17) 2-Methylphenol	8.48	107	588892	68.770	ng	# 78
18) Hexachloroethane	9.08	117	475424	71.011	ng	# 61
19) n-Nitroso-di-n-propylamine	8.85	70	669181	72.046	ng	# 84
20) 3+4-Methylphenols	8.82	107	832261	70.723	ng	# 83
22) Acetophenone	8.87	105	1252107	74.434	ng	# 95
24) Nitrobenzene	9.26	77	1088522	74.980	ng	94
25) Isophorone	9.78	82	1522751	75.808	ng	# 95
26) 2-Nitrophenol	9.96	139	487651	79.851	ng	# 55
27) 2,4-Dimethylphenol	10.01	122	618211	75.723	ng	# 79
28) bis(2-Chloroethoxy)methane	10.25	93	901404	77.577	ng	# 97
29) 2,4-Dichlorophenol	10.49	162	1008713	91.395	ng	93
30) 1,2,4-Trichlorobenzene	10.69	180	1520598	116.507	ng	95
31) Naphthalene	10.88	128	2386287	75.016	ng	97
32) Benzoic acid	10.26	122	538227m	81.166	ng	
33) 4-Chloroaniline	11.01	127	1034616	75.888	ng	# 89
34) Hexachlorobutadiene	11.13	225	1381985	122.841	ng	99
35) Caprolactam	11.87	113	225948m	79.855	ng	
36) 4-Chloro-3-methylphenol	12.14	107	901596	71.637	ng	94
37) 2-Methylnaphthalene	12.49	142	1962988	83.182	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	2038554	116.047	ng	# 100
40) Hexachlorocyclopentadiene	12.81	237	873946	160.324	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016487.D
 Acq On : 21 Aug 2018 15:46
 Operator : SJ/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:06:57 PM

Quant Time: Aug 21 16:27:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 14:43:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	1278514	116.448	nq	98
43) 2,4,5-Trichlorophenol	13.19	196	1256420	117.296	nq #	94
45) 1,1'-Biphenyl	13.50	154	3202316	80.629	nq	98
46) 2-Chloronaphthalene	13.55	162	2617536	84.478	nq	98
47) 2-Nitroaniline	13.77	65	659555	61.383	nq	85
48) Acenaphthylene	14.39	152	3780224	81.137	nq	99
49) Dimethylphthalate	14.13	163	3442347	81.804	nq #	97
50) 2,6-Dinitrotoluene	14.27	165	693816	86.868	nq #	86
51) Acenaphthene	14.73	154	2261868	80.446	nq	97
52) 3-Nitroaniline	14.61	138	594702	75.573	nq #	80
54) Dibenzofuran	15.07	168	4690948	94.686	nq	95
55) 4-Nitrophenol	14.92	139	405907	78.969	nq #	84
56) 2,4-Dinitrotoluene	15.07	165	1244588	93.222	nq #	89
57) Fluorene	15.72	166	3530463	93.275	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.30	232	1130872	112.682	nq #	100
59) Diethylphthalate	15.49	149	3572522	74.818	nq	97
60) 4-Chlorophenyl-phenylether	15.70	204	2869528	119.835	nq	91
61) 4-Nitroaniline	15.78	138	579980	76.228	nq #	1
62) Azobenzene	16.00	77	3687349	66.962	nq	90
64) 4,6-Dinitro-2-methylphenol	15.83	198	848465	96.403	nq #	63
65) n-Nitrosodiphenylamine	15.93	169	3202694	74.990	nq	99
66) 4-Bromophenyl-phenylether	16.59	248	1694202	99.804	nq	97
67) Hexachlorobenzene	16.71	284	2021802	121.368	nq	95
68) Atrazine	16.87	200	1106674	70.686	nq	90
69) Pentachlorophenol	17.06	266	1018578	106.788	nq	99
70) Phenanthrene	17.45	178	5847570	81.776	nq	99
71) Anthracene	17.54	178	5912839	83.296	nq	99
72) Carbazole	17.82	167	5174152	75.308	nq	98
73) Di-n-butylphthalate	18.35	149	6333775	64.235	nq #	97
74) Fluoranthene	19.44	202	8655024	100.195	nq	94
76) Benzidine	19.64	184	2670704	62.128	nq	99
77) Pyrene	19.81	202	9183756	71.253	nq	99
79) Butylbenzylphthalate	20.67	149	3199240	52.961	nq #	84
80) Benzo(a)anthracene	21.53	228	9668641	76.512	nq	98
81) 3,3'-Dichlorobenzidine	21.47	252	4160311	92.478	nq #	97
82) Chrysene	21.59	228	9088492	76.530	nq	98
83) Bis(2-ethylhexyl)phthalate	21.43	149	4868021	56.492	nq #	98
84) Di-n-octyl phthalate	22.32	149	7941777	61.445	nq #	92
85) Indeno(1,2,3-cd)pyrene	26.18	276	12679917	130.534	nq #	93
87) Benzo(b)fluoranthene	23.15	252	10130590	79.146	nq #	91
88) Benzo(k)fluoranthene	23.20	252	9828276	76.833	nq #	93
89) Benzo(a)pyrene	23.74	252	9493832	81.941	nq #	94
90) Dibenzo(a,h)anthracene	26.17	278	10911462	105.020	nq #	89
91) Benzo(g,h,i)perylene	26.89	276	9067789	97.105	nq #	81

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

