

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM083019\
 Data File : BM022444.D
 Acq On : 30 Aug 2019 17:02
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 8/31/2019 2:43:52 PM

Quant Time: Aug 30 18:25:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 16:51:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	22180	20.00	ng	0.00
21) Naphthalene-d8	10.29	136	100059	20.00	ng	0.00
39) Acenaphthene-d10	14.19	164	73161	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	175917	20.00	ng	0.00
76) Chrysene-d12	21.17	240	248399	20.00	ng	0.00
87) Perylene-d12	23.36	264	215977	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.15	112	104332	82.57	ng	0.00
7) Phenol-d6	6.72	99	147637	87.37	ng	0.00
23) Nitrobenzene-d5	8.70	82	190324	83.31	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	120674	80.61	ng	0.00
45) 2-Fluorobiphenyl	12.79	172	508117	78.68	ng	0.00
79) Terphenyl-d14	19.60	244	1453976	77.25	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.13	88	22742	39.003	ng	97
3) Pyridine	3.53	79	58900	41.995	ng	95
4) n-Nitrosodimethylamine	3.46	42	40994	42.394	ng	# 99
6) Aniline	6.88	93	92749	43.070	ng	100
8) 2-Chlorophenol	7.09	128	61303	42.149	ng	97
9) Benzaldehyde	6.70	77	48760	46.317	ng	94
10) Phenol	6.75	94	72438	42.673	ng	98
11) bis(2-Chloroethyl)ether	6.97	93	53956	42.118	ng	99
12) 1,3-Dichlorobenzene	7.40	146	72370	41.010	ng	95
13) 1,4-Dichlorobenzene	7.56	146	73442	41.074	ng	97
14) 1,2-Dichlorobenzene	7.86	146	74279	41.438	ng	98
15) Benzyl Alcohol	7.79	79	72829	45.431	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.06	45	73677	44.379	ng	96
17) 2-Methylphenol	7.98	107	55314	44.807	ng	89
18) Hexachloroethane	8.56	117	33944	40.944	ng	96
19) n-Nitroso-di-n-propylamine	8.34	70	58332	44.480	ng	97
20) 3+4-Methylphenols	8.31	107	75507	45.041	ng	99
22) Acetophenone	8.36	105	115558	41.823	ng	# 98
24) Nitrobenzene	8.74	77	91331	41.530	ng	98
25) Isophorone	9.26	82	157689	43.961	ng	99
26) 2-Nitrophenol	9.43	139	38511	43.717	ng	97
27) 2,4-Dimethylphenol	9.50	122	54791	41.697	ng	94
28) bis(2-Chloroethoxy)methane	9.73	93	86601	41.725	ng	98
29) 2,4-Dichlorophenol	9.96	162	72175	41.980	ng	96
30) 1,2,4-Trichlorobenzene	10.16	180	88923	40.617	ng	97
31) Naphthalene	10.35	128	216199	41.541	ng	100
32) Benzoic acid	9.72	122	47638	46.027	ng	96
33) 4-Chloroaniline	10.49	127	96690	43.178	ng	99
34) Hexachlorobutadiene	10.59	225	86084	39.865	ng	98
35) Caprolactam	11.34	113	21436m	49.054	ng	
36) 4-Chloro-3-methylphenol	11.62	107	77787	42.927	ng	97
37) 2-Methylnaphthalene	11.97	142	163529	41.348	ng	100
38) 1-Methylnaphthalene	12.19	142	156848	41.089	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.34	216	137617	39.635	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.29	237	52972	43.243	ng	97
43) 2,4,6-Trichlorophenol	12.60	196	78467	41.868	ng	98
44) 2,4,5-Trichlorophenol	12.69	196	76782	41.317	ng	98
46) 1,1'-Biphenyl	13.00	154	240517	40.234	ng	99
47) 2-Chloronaphthalene	13.05	162	190074	40.044	ng	98
48) 2-Nitroaniline	13.30	65	68176	44.596	ng	98
49) Acenaphthylene	13.90	152	291024	40.624	ng	99
50) Dimethylphthalate	13.66	163	268611	42.213	ng	99
51) 2,6-Dinitrotoluene	13.80	165	53082	42.466	ng	96
52) Acenaphthene	14.25	154	174694	40.660	ng	98
53) 3-Nitroaniline	14.14	138	53355	44.455	ng	99
54) 2,4-Dinitrophenol	14.38	184	28412	42.174	ng	95
55) Dibenzofuran	14.59	168	295163	40.781	ng	100
56) 4-Nitrophenol	14.48	139	33652	44.635	ng	98
57) 2,4-Dinitrotoluene	14.61	165	77464	43.105	ng	# 94
58) Fluorene	15.24	166	265065	40.992	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.83	232	85111	40.639	ng	96
60) Diethylphthalate	15.03	149	282205	42.459	ng	97
61) 4-Chlorophenyl-phenylether	15.24	204	179167	40.294	ng	100
62) 4-Nitroaniline	15.33	138	52973	47.156	ng	94
63) Azobenzene	15.54	77	279269	42.092	ng	99
65) 4,6-Dinitro-2-methylphenol	15.38	198	45090	42.954	ng	95
66) n-Nitrosodiphenylamine	15.47	169	210715	39.674	ng	99
67) 4-Bromophenyl-phenylether	16.14	248	110715	40.000	ng	95
68) Hexachlorobenzene	16.24	284	123657	39.441	ng	98
69) Atrazine	16.44	200	87934	38.887	ng	99
70) Pentachlorophenol	16.61	266	60399	43.070	ng	98
71) Phenanthrene	17.00	178	416342	40.602	ng	99
72) Anthracene	17.09	178	414688	41.103	ng	99
73) Carbazole	17.38	167	353516	42.391	ng	99
74) Di-n-butylphthalate	17.92	149	503055	44.487	ng	100
75) Fluoranthene	19.03	202	604308	44.254	ng	98
77) Benzidine	19.24	184	201541	48.184	ng	98
78) Pyrene	19.39	202	624403	36.885	ng	99
80) Butylbenzylphthalate	20.30	149	251161	42.513	ng	97
81) Benzo(a)anthracene	21.16	228	741861	41.609	ng	99
82) 3,3'-Dichlorobenzidine	21.10	252	247198	44.697	ng	97
83) Chrysene	21.21	228	707401	41.685	ng	100
84) Bis(2-ethylhexyl)phthalate	21.07	149	370680	46.209	ng	98
85) Di-n-octyl phthalate	21.94	149	614504	47.698	ng	99
86) Indeno(1,2,3-cd)pyrene	25.54	276	595600	34.892	ng	99
88) Benzo(b)fluoranthene	22.70	252	646697	43.101	ng	99
89) Benzo(k)fluoranthene	22.75	252	643699	43.247	ng	99
90) Benzo(a)pyrene	23.27	252	560431	42.082	ng	98
91) Dibenzo(a,h)anthracene	25.54	278	496028	37.352	ng	99
92) Benzo(g,h,i)perylene	26.21	276	412975	35.682	ng	98

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

