

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081719\
 Data File : BG042281.D
 Acq On : 17 Aug 2019 9:03
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Jagrut
 8/19/2019 2:19:40 PM

Quant Time: Aug 19 06:35:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 08:57:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	54197	20.00	ng	-0.04
21) Naphthalene-d8	10.83	136	228724	20.00	ng	-0.04
39) Acenaphthene-d10	14.67	164	171130	20.00	ng	-0.04
64) Phenanthrene-d10	17.42	188	397125	20.00	ng	-0.04
76) Chrysene-d12	21.70	240	409568	20.00	ng	-0.04
87) Perylene-d12	24.94	264	493696	20.00	ng	-0.07

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.56	112	232035	83.30	ng	-0.03
7) Phenol-d6	7.17	99	308489	82.14	ng	-0.03
23) Nitrobenzene-d5	9.17	82	272093	81.86	ng	-0.04
42) 2,4,6-Tribromophenol	16.16	330	192236	98.90	ng	-0.03
45) 2-Fluorobiphenyl	13.29	172	842935	86.88	ng	-0.04
79) Terphenyl-d14	20.03	244	1520431	84.95	ng	-0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.43	88	48542	37.024	ng	# 78
3) Pyridine	3.83	79	125949	35.031	ng	# 76
4) n-Nitrosodimethylamine	3.75	42	41189	28.899	ng	# 63
6) Aniline	7.32	93	193033	36.504	ng	91
8) 2-Chlorophenol	7.57	128	139545	43.748	ng	86
9) Benzaldehyde	7.14	77	83290	31.520	ng	86
10) Phenol	7.19	94	158884	40.667	ng	93
11) bis(2-Chloroethyl)ether	7.42	93	124256	38.686	ng	88
12) 1,3-Dichlorobenzene	7.90	146	170794m	41.026	ng	
13) 1,4-Dichlorobenzene	8.05	146	173655	41.687	ng	94
14) 1,2-Dichlorobenzene	8.37	146	165805	41.713	ng	92
15) Benzyl Alcohol	8.25	79	103203	34.931	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.54	45	163847	28.673	ng	89
17) 2-Methylphenol	8.45	107	112654	39.648	ng	97
18) Hexachloroethane	9.10	117	58686	41.140	ng	89
19) n-Nitroso-di-n-propylamine	8.82	70	98063	35.967	ng	# 80
20) 3+4-Methylphenols	8.79	107	159764	41.089	ng	97
22) Acetophenone	8.83	105	201971	39.478	ng	# 87
24) Nitrobenzene	9.22	77	137606	39.299	ng	89
25) Isophorone	9.74	82	271620	37.567	ng	96
26) 2-Nitrophenol	9.94	139	87525	53.858	ng	# 87
27) 2,4-Dimethylphenol	10.00	122	115307	40.202	ng	98
28) bis(2-Chloroethoxy)methane	10.23	93	173428	38.384	ng	99
29) 2,4-Dichlorophenol	10.48	162	158767	45.455	ng	96
30) 1,2,4-Trichlorobenzene	10.70	180	186786	42.171	ng	98
31) Naphthalene	10.88	128	447148	42.716	ng	96
32) Benzoic acid	10.13	122	65472	34.871	ng	95
33) 4-Chloroaniline	10.98	127	201183	40.526	ng	# 92
34) Hexachlorobutadiene	11.17	225	123368	41.269	ng	98
35) Caprolactam	11.76	113	53969	44.461	ng	# 51
36) 4-Chloro-3-methylphenol	12.12	107	152145	43.936	ng	94
37) 2-Methylnaphthalene	12.49	142	359638	43.404	ng	97
38) 1-Methylnaphthalene	12.71	142	344838	43.916	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	220820	40.764	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.84	237	103505	33.985	nq	99
43) 2,4,6-Trichlorophenol	13.10	196	139586	40.770	nq	98
44) 2,4,5-Trichlorophenol	13.18	196	151148	42.482	nq #	92
46) 1,1'-Biphenyl	13.50	154	460668	41.876	nq	95
47) 2-Chloronaphthalene	13.55	162	388370	41.466	nq	95
48) 2-Nitroaniline	13.74	65	94477	39.773	nq #	73
49) Acenaphthylene	14.39	152	613411	43.784	nq	97
50) Dimethylphthalate	14.12	163	514374	44.011	nq	99
51) 2,6-Dinitrotoluene	14.23	165	122524	50.135	nq #	79
52) Acenaphthene	14.73	154	365332	40.094	nq	98
53) 3-Nitroaniline	14.57	138	119993	45.895	nq #	80
54) 2,4-Dinitrophenol	14.78	184	52227	43.192	nq #	80
55) Dibenzofuran	15.07	168	611977	43.909	nq	92
56) 4-Nitrophenol	14.88	139	88533	44.610	nq #	84
57) 2,4-Dinitrotoluene	15.03	165	175066	52.090	nq #	83
58) Fluorene	15.72	166	499469	44.656	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.30	232	140881	39.980	nq #	91
60) Diethylphthalate	15.48	149	501572	43.713	nq	99
61) 4-Chlorophenyl-phenylether	15.71	204	291744	42.727	nq	94
62) 4-Nitroaniline	15.73	138	128495	45.375	nq	92
63) Azobenzene	16.00	77	349097	37.244	nq	87
65) 4,6-Dinitro-2-methylphenol	15.80	198	92143	47.491	nq	80
66) n-Nitrosodiphenylamine	15.92	169	460163	42.539	nq	96
67) 4-Bromophenyl-phenylether	16.61	248	203377	41.575	nq	94
68) Hexachlorobenzene	16.73	284	217342	42.469	nq #	90
69) Atrazine	16.87	200	39827	9.371	nq	100
70) Pentachlorophenol	17.08	266	101492	34.910	nq	99
71) Phenanthrene	17.47	178	840133	44.135	nq	94
72) Anthracene	17.55	178	845466	44.534	nq	94
73) Carbazole	17.82	167	763856	44.829	nq #	93
74) Di-n-butylphthalate	18.38	149	894750	46.318	nq #	94
75) Fluoranthene	19.47	202	1060532	45.835	nq	90
77) Benzidine	19.65	184	464804	34.769	nq	98
78) Pyrene	19.84	202	1059719	43.670	nq	91
80) Butylbenzylphthalate	20.72	149	419277	45.065	nq #	84
81) Benzo(a)anthracene	21.68	228	1062957	43.866	nq	93
82) 3,3'-Dichlorobenzidine	21.59	252	410337	42.177	nq	96
83) Chrysene	21.75	228	1026301	44.184	nq	92
84) Bis(2-ethylhexyl)phthalate	21.58	149	582480	45.387	nq	96
85) Di-n-octyl phthalate	22.81	149	992720	45.963	nq #	86
86) Indeno(1,2,3-cd)pyrene	28.66	276	1334936	43.415	nq #	94
88) Benzo(b)fluoranthene	23.92	252	1149768	42.572	nq #	96
89) Benzo(k)fluoranthene	23.99	252	1084532	41.225	nq #	96
90) Benzo(a)pyrene	24.80	252	1084354	41.864	nq #	95
91) Dibenzo(a,h)anthracene	28.73	278	1078945	42.422	nq #	96
92) Benzo(g,h,i)perylene	29.82	276	1086158	42.745	nq #	93

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

