

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011420\
 Data File : BP001590.D
 Acq On : 14 Jan 2020 16:35
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/16/2020 1:45:45 PM

Quant Time: Jan 14 21:54:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 21:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	25770	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	100002	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	76414	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	188137	20.00	ng	0.00
76) Chrysene-d12	21.16	240	179965	20.00	ng	0.00
87) Perylene-d12	23.39	264	173274	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	107556	81.76	ng	0.00
7) Phenol-d6	6.83	99	148886	70.81	ng	0.00
23) Nitrobenzene-d5	8.79	82	179086	78.25	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	80817	69.48	ng	0.00
45) 2-Fluorobiphenyl	12.90	172	410114	78.91	ng	0.00
79) Terphenyl-d14	19.64	244	751585	76.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.23	88	21993	52.384	ng	99
3) Pyridine	3.61	79	64273	42.366	ng	97
4) n-Nitrosodimethylamine	3.53	42	51385	47.789	ng	96
6) Aniline	6.98	93	95530	36.339	ng	97
8) 2-Chlorophenol	7.21	128	56685	36.908	ng	96
9) Benzaldehyde	6.79	77	42286	33.672	ng	94
10) Phenol	6.86	94	75762	34.835	ng	100
11) bis(2-Chloroethyl)ether	7.08	93	64304	37.698	ng	94
12) 1,3-Dichlorobenzene	7.53	146	76111	39.424	ng	94
13) 1,4-Dichlorobenzene	7.68	146	75580	37.861	ng	95
14) 1,2-Dichlorobenzene	7.99	146	71643	37.137	ng	98
15) Benzyl Alcohol	7.88	79	63583	32.113	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.18	45	124800	39.384	ng	96
17) 2-Methylphenol	8.09	107	52488	34.759	ng	99
18) Hexachloroethane	8.71	117	31373	40.257	ng	96
19) n-Nitroso-di-n-propylamine	8.45	70	55201	34.438	ng	95
20) 3+4-Methylphenols	8.42	107	71141	33.681	ng	97
22) Acetophenone	8.46	105	103693	36.990	ng	# 94
24) Nitrobenzene	8.83	77	87207	37.667	ng	95
25) Isophorone	9.36	82	145417	35.434	ng	99
26) 2-Nitrophenol	9.53	139	31480	35.979	ng	98
27) 2,4-Dimethylphenol	9.61	122	46031	35.215	ng	96
28) bis(2-Chloroethoxy)methane	9.85	93	87728	36.491	ng	97
29) 2,4-Dichlorophenol	10.07	162	63690	35.597	ng	96
30) 1,2,4-Trichlorobenzene	10.28	180	83690	38.746	ng	94
31) Naphthalene	10.46	128	195067	36.960	ng	99
32) Benzoic acid	9.75	122	29231m	27.117	ng	
33) 4-Chloroaniline	10.58	127	78610	32.349	ng	97
34) Hexachlorobutadiene	10.76	225	60892	39.825	ng	96
35) Caprolactam	11.36	113	18026	27.744	ng	93
36) 4-Chloro-3-methylphenol	11.72	107	69976	32.588	ng	98
37) 2-Methylnaphthalene	12.09	142	149102	35.448	ng	94
38) 1-Methylnaphthalene	12.31	142	141678	35.003	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.46	216	104122	39.778	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.45	237	56799	42.777	nq	99
43) 2,4,6-Trichlorophenol	12.71	196	61272	36.294	nq	95
44) 2,4,5-Trichlorophenol	12.78	196	65959	36.767	nq	94
46) 1,1'-Biphenyl	13.12	154	209288	38.221	nq	99
47) 2-Chloronaphthalene	13.15	162	175795	38.628	nq	98
48) 2-Nitroaniline	13.36	65	64472	36.059	nq	97
49) Acenaphthylene	14.00	152	262435	37.391	nq	99
50) Dimethylphthalate	13.76	163	220301	35.003	nq	98
51) 2,6-Dinitrotoluene	13.87	165	48275	36.170	nq	98
52) Acenaphthene	14.35	154	160649	36.958	nq	99
53) 3-Nitroaniline	14.20	138	45254	32.114	nq	92
54) 2,4-Dinitrophenol	14.40	184	20722	27.613	nq	# 79
55) Dibenzofuran	14.69	168	270427	36.789	nq	97
56) 4-Nitrophenol	14.52	139	32499	28.881	nq	97
57) 2,4-Dinitrotoluene	14.66	165	66776	34.194	nq	# 99
58) Fluorene	15.35	166	220424	37.058	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.93	232	63097	34.061	nq	99
60) Diethylphthalate	15.14	149	226239	34.949	nq	100
61) 4-Chlorophenyl-phenylether	15.35	204	127061	37.325	nq	98
62) 4-Nitroaniline	15.37	138	49912	31.066	nq	97
63) Azobenzene	15.65	77	248996	38.496	nq	97
65) 4,6-Dinitro-2-methylphenol	15.43	198	34507	33.925	nq	92
66) n-Nitrosodiphenylamine	15.56	169	188753	37.870	nq	99
67) 4-Bromophenyl-phenylether	16.25	248	81099	38.211	nq	96
68) Hexachlorobenzene	16.36	284	94827	38.455	nq	96
69) Atrazine	16.53	200	64657	35.322	nq	97
70) Pentachlorophenol	16.71	266	45835	33.657	nq	97
71) Phenanthrene	17.09	178	382414	37.376	nq	99
72) Anthracene	17.19	178	370139	37.161	nq	100
73) Carbazole	17.46	167	325672	34.827	nq	99
74) Di-n-butylphthalate	18.05	149	402699	36.399	nq	100
75) Fluoranthene	19.07	202	478935	35.575	nq	100
77) Benzidine	19.26	184	144701	35.214	nq	96
78) Pyrene	19.43	202	493117	37.301	nq	99
80) Butylbenzylphthalate	20.33	149	184830	36.786	nq	98
81) Benzo(a)anthracene	21.14	228	457762	36.523	nq	99
82) 3,3'-Dichlorobenzidine	21.08	252	159225	33.878	nq	97
83) Chrysene	21.19	228	453213	37.105	nq	99
84) Bis(2-ethylhexyl)phthalate	21.10	149	262114	38.054	nq	97
85) Di-n-octyl phthalate	21.98	149	409139	37.253	nq	96
86) Indeno(1,2,3-cd)pyrene	25.66	276	436596	30.536	nq	98
88) Benzo(b)fluoranthene	22.72	252	431291	39.500	nq	99
89) Benzo(k)fluoranthene	22.77	252	433033	40.676	nq	99
90) Benzo(a)pyrene	23.30	252	393853	37.990	nq	98
91) Dibenzo(a,h)anthracene	25.67	278	364373	34.065	nq	# 97
92) Benzo(g,h,i)perylene	26.34	276	363511	34.185	nq	98

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

