

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020521\
 Data File : BP004677.D
 Acq On : 05 Feb 2021 13:35
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC005

Quant Time: Feb 05 16:04:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 16:02:44 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	43470	20.00	ng	# 0.00
21) Naphthalene-d8	10.569	136	169798	20.00	ng	# 0.00
39) Acenaphthene-d10	14.422	164	110896	20.00	ng	0.00
64) Phenanthrene-d10	17.181	188	246299	20.00	ng	0.00
76) Chrysene-d12	21.257	240	271584	20.00	ng	0.00
86) Perylene-d12	23.569	264	269522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	32177	12.90	ng	0.00
7) Phenol-d6	6.940	99	42861	12.55	ng	0.00
23) Nitrobenzene-d5	8.934	82	39569	13.17	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	14088	7.31	ng	0.00
45) 2-Fluorobiphenyl	13.046	172	97768	11.46	ng	0.00
79) Terphenyl-d14	19.739	244	172644	11.61	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	7244	7.18	ng	# 87
3) Pyridine	3.705	79	14538	5.38	ng	89
4) n-Nitrosodimethylamine	3.605	42	6521	7.63	ng	# 62
6) Aniline	7.105	93	23735	6.11	ng	# 84
8) 2-Chlorophenol	7.340	128	17054	6.02	ng	89
10) Phenol	6.969	94	21654	6.61	ng	82
11) bis(2-Chloroethyl)ether	7.205	93	16686	6.37	ng	97
12) 1,3-Dichlorobenzene	7.664	146	21797	6.22	ng	# 92
13) 1,4-Dichlorobenzene	7.811	146	21636	6.11	ng	94
14) 1,2-Dichlorobenzene	8.128	146	21126	6.24	ng	88
15) Benzyl Alcohol	8.016	79	14711	6.65	ng	# 86
16) 2,2'-oxybis(1-Chloropr...	8.316	45	13188	4.74	ng	85
17) 2-Methylphenol	8.216	107	14057	6.08	ng	# 89
18) Hexachloroethane	8.852	117	7486	6.11	ng	92
19) n-Nitroso-di-n-propyla...	8.581	70	12060	6.29	ng	# 81
20) 3+4-Methylphenols	8.540	107	17601	5.58	ng	89
22) Acetophenone	8.599	105	25752	6.14	ng	# 88
24) Nitrobenzene	8.975	77	20374	6.93	ng	# 79
25) Isophorone	9.499	82	30936	5.71	ng	# 87
26) 2-Nitrophenol	9.687	139	7106	4.39	ng	# 72
27) 2,4-Dimethylphenol	9.746	122	12772	5.69	ng	# 83
28) bis(2-Chloroethoxy)met...	9.987	93	21466	5.78	ng	100
29) 2,4-Dichlorophenol	10.216	162	14376	4.83	ng	88
30) 1,2,4-Trichlorobenzene	10.434	180	19789	5.52	ng	95
31) Naphthalene	10.622	128	55857	5.99	ng	98
33) 4-Chloroaniline	10.734	127	20970	5.21	ng	# 85
34) Hexachlorobutadiene	10.910	225	12664	5.26	ng	95
35) Caprolactam	11.487	113	3461	3.64	ng	# 69
36) 4-Chloro-3-methylphenol	11.857	107	17260	6.08	ng	# 81
37) 2-Methylnaphthalene	12.240	142	38204	5.57	ng	# 93
38) 1-Methylnaphthalene	12.457	142	36346	5.58	ng	# 94
40) 1,2,4,5-Tetrachloroben...	12.604	216	21793	5.47	ng	# 97
41) Hexachlorocyclopentadiene	12.587	237	11931	8.39	ng	93
43) 2,4,6-Trichlorophenol	12.851	196	12398	4.78	ng	92
44) 2,4,5-Trichlorophenol	12.916	196	13736	5.10	ng	# 88
46) 1,1'-Biphenyl	13.257	154	50526	5.70	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020521\
 Data File : BP004677.D
 Acq On : 05 Feb 2021 13:35
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC005

Quant Time: Feb 05 16:04:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 16:02:44 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.293	162	41415	5.64	ng	96
48) 2-Nitroaniline	13.498	65	9451	5.57	ng #	69
49) Acenaphthylene	14.146	152	58443	5.33	ng	99
50) Dimethylphthalate	13.881	163	52577	5.72	ng	99
51) 2,6-Dinitrotoluene	13.998	165	9545	4.87	ng #	65
52) Acenaphthene	14.487	154	38891	5.82	ng	91
53) 3-Nitroaniline	14.328	138	9226	4.33	ng #	65
54) 2,4-Dinitrophenol	14.528	184	4751	5.37	ng #	71
55) Dibenzofuran	14.822	168	61708	5.70	ng	96
56) 4-Nitrophenol	14.628	139	7006	4.73	ng #	55
57) 2,4-Dinitrotoluene	14.781	165	12571	4.69	ng #	60
58) Fluorene	15.475	166	49259	5.70	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.045	232	12738	5.04	ng	95
60) Diethylphthalate	15.251	149	51358	5.71	ng	99
61) 4-Chlorophenyl-phenyle...	15.475	204	28202	5.78	ng	97
62) 4-Nitroaniline	15.487	138	9941	4.41	ng #	60
63) Azobenzene	15.763	77	47338	6.96	ng	83
65) 4,6-Dinitro-2-methylph...	15.545	198	6450	4.50	ng	89
66) n-Nitrosodiphenylamine	15.687	169	41636	5.45	ng	97
67) 4-Bromophenyl-phenylether	16.369	248	17546	5.33	ng	90
68) Hexachlorobenzene	16.481	284	18367	4.67	ng #	90
69) Atrazine	16.640	200	13356	4.97	ng	96
70) Pentachlorophenol	16.822	266	8624	4.89	ng	95
71) Phenanthrene	17.216	178	82486	5.84	ng	99
72) Anthracene	17.310	178	76403	5.58	ng	98
73) Carbazole	17.581	167	70138	5.50	ng	97
74) Di-n-butylphthalate	18.145	149	78175	5.07	ng #	96
75) Fluoranthene	19.186	202	96273	5.60	ng	97
77) Benzidine	19.369	184	20735	2.92	ng	96
78) Pyrene	19.539	202	99350	5.38	ng	99
80) Butylbenzylphthalate	20.422	149	34450	4.61	ng #	80
81) Benzo(a)anthracene	21.245	228	103745	5.62	ng	97
82) 3,3'-Dichlorobenzidine	21.180	252	31380	4.71	ng	99
83) Chrysene	21.298	228	101885	5.79	ng	98
84) Bis(2-ethylhexyl)phtha...	21.180	149	54082	5.07	ng	97
85) Di-n-octyl phthalate	22.075	149	86568	4.78	ng #	91
87) Indeno(1,2,3-cd)pyrene	25.933	276	112800	5.42	ng	96
88) Benzo(b)fluoranthene	22.869	252	107060	6.14	ng #	98
89) Benzo(k)fluoranthene	22.916	252	100483	6.04	ng #	97
90) Benzo(a)pyrene	23.463	252	93341	5.74	ng	98
91) Dibenzo(a,h)anthracene	25.957	278	95821	5.58	ng #	95
92) Benzo(g,h,i)perylene	26.663	276	92102	5.45	ng #	94

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

