

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072221\
 Data File : BP006274.D
 Acq On : 22 Jul 2021 11:03
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 7/23/2021 11:51:39 AM

Quant Time: Jul 22 12:50:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 12:49:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	248161	20.00	ng	0.00
21) Naphthalene-d8	10.587	136	1050366	20.00	ng	# 0.00
39) Acenaphthene-d10	14.428	164	631579	20.00	ng	0.00
64) Phenanthrene-d10	17.175	188	1259407	20.00	ng	# 0.00
76) Chrysene-d12	21.274	240	1239344	20.00	ng	# 0.00
86) Perylene-d12	23.627	264	1250034	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	318181	21.36	ng	0.00
7) Phenol-d6	6.969	99	448528	22.02	ng	0.00
23) Nitrobenzene-d5	8.952	82	393105	22.54	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	100605	15.48	ng	0.00
45) 2-Fluorobiphenyl	13.051	172	873189	21.04	ng	0.00
79) Terphenyl-d14	19.751	244	1304179	21.05	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.340	88	68000	10.92	ng	93
3) Pyridine	3.734	79	193985	11.19	ng	97
4) n-Nitrosodimethylamine	3.646	42	65895	11.21	ng	# 75
6) Aniline	7.128	93	242673	10.76	ng	96
8) 2-Chlorophenol	7.363	128	176134	10.71	ng	95
9) Benzaldehyde	6.946	77	142258m	12.13	ng	
10) Phenol	6.993	94	219947	10.89	ng	96
11) bis(2-Chloroethyl)ether	7.234	93	175772	11.23	ng	93
12) 1,3-Dichlorobenzene	7.687	146	191409	10.62	ng	99
13) 1,4-Dichlorobenzene	7.834	146	195555	10.71	ng	97
14) 1,2-Dichlorobenzene	8.146	146	188246	10.76	ng	98
15) Benzyl Alcohol	8.040	79	159690	10.89	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.334	45	192995	12.24	ng	94
17) 2-Methylphenol	8.240	107	144200	10.76	ng	97
18) Hexachloroethane	8.869	117	70153	10.80	ng	# 89
19) n-Nitroso-di-n-propyla...	8.604	70	141106	11.31	ng	# 96
20) 3+4-Methylphenols	8.563	107	194513	10.78	ng	95
22) Acetophenone	8.622	105	264236	10.42	ng	# 98
24) Nitrobenzene	8.993	77	207049	11.16	ng	93
25) Isophorone	9.522	82	363957	10.23	ng	94
26) 2-Nitrophenol	9.704	139	70730	10.15	ng	96
27) 2,4-Dimethylphenol	9.769	122	144188	10.24	ng	98
28) bis(2-Chloroethoxy)met...	10.010	93	214927	10.70	ng	98
29) 2,4-Dichlorophenol	10.234	162	142350	9.79	ng	96
30) 1,2,4-Trichlorobenzene	10.446	180	165855	10.09	ng	97
31) Naphthalene	10.634	128	577118	10.47	ng	99
32) Benzoic acid	9.840	122	83256	9.53	ng	100
33) 4-Chloroaniline	10.746	127	222054	9.89	ng	96
34) Hexachlorobutadiene	10.922	225	96659	9.92	ng	97
35) Caprolactam	11.522	113	43873	10.03	ng	83
36) 4-Chloro-3-methylphenol	11.869	107	175367	10.37	ng	95
37) 2-Methylnaphthalene	12.245	142	397751	10.33	ng	99
38) 1-Methylnaphthalene	12.463	142	374028	10.27	ng	98
40) 1,2,4,5-Tetrachloroben...	12.610	216	170518	9.97	ng	# 97
41) Hexachlorocyclopentadiene	12.593	237	88269	9.42	ng	98
43) 2,4,6-Trichlorophenol	12.857	196	107249	10.00	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.922	196	120296	10.43	ng	# 95
46) 1,1'-Biphenyl	13.263	154	484429	10.41	ng	98
47) 2-Chloronaphthalene	13.298	162	368982	10.29	ng	95
48) 2-Nitroaniline	13.504	65	92913	11.03	ng	90
49) Acenaphthylene	14.145	152	586124	10.27	ng	100
50) Dimethylphthalate	13.892	163	459384	10.37	ng	99
51) 2,6-Dinitrotoluene	14.004	165	91587	10.53	ng	# 84
52) Acenaphthene	14.487	154	367947	10.43	ng	97
53) 3-Nitroaniline	14.334	138	95193	10.33	ng	# 88
54) 2,4-Dinitrophenol	14.534	184	38097	10.33	ng	# 89
55) Dibenzofuran	14.822	168	580038	10.42	ng	97
56) 4-Nitrophenol	14.634	139	79028	9.67	ng	86
57) 2,4-Dinitrotoluene	14.792	165	113130	10.08	ng	91
58) Fluorene	15.475	166	461992	10.36	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	100739	10.09	ng	# 78
60) Diethylphthalate	15.263	149	468970	10.25	ng	99
61) 4-Chlorophenyl-phenyle...	15.475	204	216568	10.26	ng	93
62) 4-Nitroaniline	15.492	138	97430	10.12	ng	84
63) Azobenzene	15.769	77	497173	10.80	ng	93
65) 4,6-Dinitro-2-methylph...	15.551	198	55507	9.47	ng	82
66) n-Nitrosodiphenylamine	15.692	169	396898	10.27	ng	99
67) 4-Bromophenyl-phenylether	16.375	248	107031	8.44	ng	94
68) Hexachlorobenzene	16.475	284	142247	9.78	ng	92
69) Atrazine	16.651	200	125351	9.93	ng	96
70) Pentachlorophenol	16.822	266	76267	9.36	ng	98
71) Phenanthrene	17.216	178	730315	10.59	ng	100
72) Anthracene	17.310	178	687366	10.22	ng	100
73) Carbazole	17.586	167	684938	10.16	ng	99
74) Di-n-butylphthalate	18.157	149	746365	9.69	ng	100
75) Fluoranthene	19.192	202	803920	10.15	ng	94
77) Benzidine	19.380	184	349453	9.69	ng	100
78) Pyrene	19.545	202	837910	10.24	ng	99
80) Butylbenzylphthalate	20.439	149	316906	9.14	ng	93
81) Benzo(a)anthracene	21.257	228	820757	10.32	ng	99
82) 3,3'-Dichlorobenzidine	21.198	252	194754	8.84	ng	# 96
83) Chrysene	21.310	228	804130	10.44	ng	100
84) Bis(2-ethylhexyl)phtha...	21.210	149	480289	9.19	ng	# 97
85) Di-n-octyl phthalate	22.121	149	784162	8.95	ng	# 96
87) Indeno(1,2,3-cd)pyrene	26.062	276	883388	9.98	ng	# 89
88) Benzo(b)fluoranthene	22.910	252	817218	10.13	ng	# 95
89) Benzo(k)fluoranthene	22.957	252	780522	10.39	ng	# 95
90) Benzo(a)pyrene	23.521	252	718006	9.92	ng	# 95
91) Dibenzo(a,h)anthracene	26.086	278	769262	10.04	ng	# 92
92) Benzo(g,h,i)perylene	26.809	276	734784	9.85	ng	# 91

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

