

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042121\
 Data File : BP005356.D
 Acq On : 21 Apr 2021 12:58
 Operator : CG/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

mohammad
 4/22/2021 12:10:57 PM

Quant Time: Apr 21 15:46:43 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 21 15:45:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.645	152	43871	20.00	ng	0.00
21) Naphthalene-d8	10.433	136	150762	20.00	ng	0.00
39) Acenaphthene-d10	14.310	164	111079	20.00	ng	0.00
64) Phenanthrene-d10	17.074	188	269974	20.00	ng	0.00
76) Chrysene-d12	21.174	240	358487	20.00	ng	0.00
86) Perylene-d12	23.439	264	382977	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.251	112	22420	8.13	ng	0.00
7) Phenol-d6	6.834	99	30920	9.71	ng	0.00
23) Nitrobenzene-d5	8.810	82	25817	4.57	ng	0.00
42) 2,4,6-Tribromophenol	15.816	330	17063	6.60	ng	0.00
45) 2-Fluorobiphenyl	12.922	172	77168	9.19	ng	0.00
79) Terphenyl-d14	19.651	244	165982	8.96	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.187	88	6396	3.74	ng	# 85
3) Pyridine	3.628	79	10431	2.89	ng	# 78
4) n-Nitrosodimethylamine	3.510	42	6532	5.36	ng	# 17
6) Aniline	6.987	93	16234	5.94	ng	# 87
8) 2-Chlorophenol	7.216	128	12398	4.67	ng	84
9) Benzaldehyde	6.804	77	10702	7.30	ng	90
10) Phenol	6.863	94	15040	4.76	ng	80
11) bis(2-Chloroethyl)ether	7.087	93	11755	5.18	ng	96
12) 1,3-Dichlorobenzene	7.534	146	15506	4.41	ng	# 87
13) 1,4-Dichlorobenzene	7.681	146	15652	4.62	ng	94
14) 1,2-Dichlorobenzene	7.998	146	15084	4.75	ng	97
15) Benzyl Alcohol	7.904	79	11379	8.41	ng	84
16) 2,2'-oxybis(1-Chloropr...	8.187	45	9807	7.30	ng	97
17) 2-Methylphenol	8.104	107	9678	6.26	ng	# 87
18) Hexachloroethane	8.716	117	5258	4.40	ng	87
19) n-Nitroso-di-n-propyla...	8.457	70	8225	6.38	ng	95
20) 3+4-Methylphenols	8.439	107	12391	6.87	ng	92
22) Acetophenone	8.481	105	19118	2.56	ng	# 93
24) Nitrobenzene	8.851	77	13791	2.54	ng	94
25) Isophorone	9.381	82	20957	3.89	ng	99
26) 2-Nitrophenol	9.569	139	4562	2.93	ng	# 87
27) 2,4-Dimethylphenol	9.634	122	8480	3.88	ng	87
28) bis(2-Chloroethoxy)met...	9.869	93	12140	4.35	ng	95
29) 2,4-Dichlorophenol	10.116	162	10993	4.05	ng	93
30) 1,2,4-Trichlorobenzene	10.298	180	16671	4.40	ng	93
31) Naphthalene	10.486	128	37898	4.28	ng	96
33) 4-Chloroaniline	10.622	127	11986	4.27	ng	97
34) Hexachlorobutadiene	10.763	225	14355	4.08	ng	89
36) 4-Chloro-3-methylphenol	11.763	107	11051	4.94	ng	94
37) 2-Methylnaphthalene	12.116	142	24801	4.58	ng	92
38) 1-Methylnaphthalene	12.333	142	24927	4.78	ng	# 90
40) 1,2,4,5-Tetrachloroben...	12.486	216	22859	5.24	ng	# 98
43) 2,4,6-Trichlorophenol	12.745	196	10173	4.46	ng	94
44) 2,4,5-Trichlorophenol	12.822	196	10629	4.20	ng	# 91
46) 1,1'-Biphenyl	13.139	154	36991	4.84	ng	99
47) 2-Chloronaphthalene	13.180	162	29924	4.65	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
48) 2-Nitroaniline	13.398	65	4810	3.37	ng	# 90
49) Acenaphthylene	14.033	152	38294	3.99	ng	97
50) Dimethylphthalate	13.774	163	32876	4.53	ng	97
51) 2,6-Dinitrotoluene	13.898	165	6129	3.29	ng	# 76
52) Acenaphthene	14.374	154	27348	4.36	ng	94
53) 3-Nitroaniline	14.245	138	5060	3.52	ng	# 81
55) Dibenzofuran	14.716	168	48508	4.06	ng	98
57) 2,4-Dinitrotoluene	14.704	165	8110	2.88	ng	# 79
58) Fluorene	15.369	166	37243	3.78	ng	95
59) 2,3,4,6-Tetrachlorophenol	14.951	232	12130	4.03	ng	# 96
60) Diethylphthalate	15.145	149	29166	3.91	ng	# 95
61) 4-Chlorophenyl-phenyle...	15.369	204	24121	4.09	ng	96
62) 4-Nitroaniline	15.416	138	3821	2.52	ng	# 53
63) Azobenzene	15.663	77	31410	3.12	ng	94
66) n-Nitrosodiphenylamine	15.586	169	31150	4.31	ng	# 93
67) 4-Bromophenyl-phenylether	16.263	248	15637	4.24	ng	94
68) Hexachlorobenzene	16.374	284	20184	4.16	ng	97
69) Atrazine	16.545	200	11486	4.04	ng	96
71) Phenanthrene	17.115	178	65037	4.05	ng	98
72) Anthracene	17.210	178	59983	3.95	ng	95
73) Carbazole	17.492	167	52735	3.86	ng	97
74) Di-n-butylphthalate	18.045	149	39395	3.37	ng	# 93
75) Fluoranthene	19.104	202	81399	3.94	ng	98
77) Benzidine	19.298	184	24994	5.51	ng	98
78) Pyrene	19.457	202	86692	4.77	ng	99
80) Butylbenzylphthalate	20.333	149	17232	4.29	ng	93
81) Benzo(a)anthracene	21.162	228	100076	4.59	ng	99
82) 3,3'-Dichlorobenzidine	21.104	252	27489	4.77	ng	99
83) Chrysene	21.215	228	108371	4.96	ng	99
84) Bis(2-ethylhexyl)phtha...	21.086	149	24863	4.02	ng	96
87) Indeno(1,2,3-cd)pyrene	25.750	276	133086	4.27	ng	99
88) Benzo(b)fluoranthene	22.750	252	109949m	4.21	ng	
89) Benzo(k)fluoranthene	22.798	252	112643m	4.53	ng	
90) Benzo(a)pyrene	23.339	252	101665	4.35	ng	98
91) Dibenzo(a,h)anthracene	25.762	278	114984	4.22	ng	99
92) Benzo(g,h,i)perylene	26.456	276	110335	4.33	ng	95

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

