

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090821\
 Data File : BM032097.D
 Acq On : 08 Sep 2021 15:38
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

mohammad
 9/9/2021 11:51:39 AM

Quant Time: Sep 08 17:07:12 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 08 17:04:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	62416	20.000	ng	0.00
21) Naphthalene-d8	10.116	136	244783	20.000	ng	# 0.00
39) Acenaphthene-d10	14.022	164	166647	20.000	ng	0.00
64) Phenanthrene-d10	16.786	188	337195	20.000	ng	0.00
76) Chrysene-d12	21.009	240	316305	20.000	ng	0.00
86) Perylene-d12	23.103	264	300074	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.004	112	142034	36.346	ng	0.00
7) Phenol-d6	6.569	99	214609	38.300	ng	0.00
23) Nitrobenzene-d5	8.516	82	221193	40.875	ng	0.00
42) 2,4,6-Tribromophenol	15.533	330	47785	26.563	ng	0.00
45) 2-Fluorobiphenyl	12.628	172	511628	41.521	ng	0.00
79) Terphenyl-d14	19.451	244	747936	38.013	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.022	88	32348	20.841	ng	# 86
3) Pyridine	3.405	79	84243	20.670	ng	90
4) n-Nitrosodimethylamine	3.328	42	44632	20.362	ng	# 58
6) Aniline	6.716	93	118840	19.846	ng	98
8) 2-Chlorophenol	6.934	128	79769	17.626	ng	97
9) Benzaldehyde	6.534	77	72221m	21.015	ng	
10) Phenol	6.593	94	108314	19.484	ng	97
11) bis(2-Chloroethyl)ether	6.816	93	83489	19.678	ng	97
12) 1,3-Dichlorobenzene	7.251	146	98739	20.568	ng	95
13) 1,4-Dichlorobenzene	7.398	146	99222	20.141	ng	97
14) 1,2-Dichlorobenzene	7.704	146	97135	20.381	ng	98
15) Benzyl Alcohol	7.616	79	82573	19.545	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.898	45	117560	20.181	ng	97
17) 2-Methylphenol	7.816	107	72379	19.234	ng	99
18) Hexachloroethane	8.410	117	36879	19.182	ng	# 88
19) n-Nitroso-di-n-propyla...	8.175	70	76666	19.309	ng	92
20) 3+4-Methylphenols	8.145	107	95670	18.405	ng	99
22) Acetophenone	8.187	105	140826	21.176	ng	96
24) Nitrobenzene	8.557	77	118055	21.460	ng	95
25) Isophorone	9.081	82	203366	21.356	ng	99
26) 2-Nitrophenol	9.257	139	21664	9.516	ng	94
27) 2,4-Dimethylphenol	9.328	122	67684	19.468	ng	93
28) bis(2-Chloroethoxy)met...	9.569	93	101488	20.267	ng	98
29) 2,4-Dichlorophenol	9.781	162	69364	17.778	ng	95
30) 1,2,4-Trichlorobenzene	9.986	180	102003	23.954	ng	96
31) Naphthalene	10.169	128	276767	20.420	ng	100
32) Benzoic acid	9.469	122	20044	7.182	ng	99
33) 4-Chloroaniline	10.304	127	105289	19.100	ng	97
34) Hexachlorobutadiene	10.439	225	68803	26.663	ng	96
35) Caprolactam	11.098	113	20207	16.519	ng	99
36) 4-Chloro-3-methylphenol	11.433	107	85441	18.715	ng	94
37) 2-Methylnaphthalene	11.792	142	198751	20.229	ng	98
38) 1-Methylnaphthalene	12.016	142	188281	20.080	ng	99
40) 1,2,4,5-Tetrachloroben...	12.175	216	111479	23.171	ng	# 96
41) Hexachlorocyclopentadiene	12.139	237	28566	12.815	ng	100
43) 2,4,6-Trichlorophenol	12.433	196	41889	12.242	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.504	196	53513	14.397	ng	93
46) 1,1'-Biphenyl	12.839	154	255266	19.560	ng	99
47) 2-Chloronaphthalene	12.880	162	205625	19.991	ng	96
48) 2-Nitroaniline	13.110	65	48207	14.664	ng	97
49) Acenaphthylene	13.739	152	310095	19.082	ng	99
50) Dimethylphthalate	13.504	163	243570	18.915	ng	99
51) 2,6-Dinitrotoluene	13.627	165	48153	16.668	ng	85
52) Acenaphthene	14.086	154	191688	19.240	ng	96
53) 3-Nitroaniline	13.957	138	41263	14.524	ng #	91
54) 2,4-Dinitrophenol	14.180	184	11394	7.067	ng #	75
55) Dibenzofuran	14.427	168	323702	19.984	ng	94
56) 4-Nitrophenol	14.286	139	20795	9.145	ng	92
57) 2,4-Dinitrotoluene	14.427	165	58189	14.829	ng #	93
58) Fluorene	15.086	166	253578	19.346	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.669	232	40055	12.728	ng #	82
60) Diethylphthalate	14.886	149	236793	17.565	ng	96
61) 4-Chlorophenyl-phenyle...	15.092	204	136402	21.588	ng	90
62) 4-Nitroaniline	15.133	138	36660	13.187	ng	99
63) Azobenzene	15.386	77	281971	18.964	ng	92
65) 4,6-Dinitro-2-methylph...	15.198	198	19429	8.638	ng	85
66) n-Nitrosodiphenylamine	15.316	169	217647	19.168	ng	94
67) 4-Bromophenyl-phenylether	15.986	248	77600	21.625	ng	94
68) Hexachlorobenzene	16.086	284	88905	23.182	ng #	86
69) Atrazine	16.286	200	70034	19.750	ng	94
70) Pentachlorophenol	16.445	266	29848	12.449	ng	94
71) Phenanthrene	16.827	178	388510	19.398	ng	99
72) Anthracene	16.921	178	379752	19.452	ng	99
73) Carbazole	17.204	167	356744	19.016	ng	98
74) Di-n-butylphthalate	17.780	149	364330	16.174	ng	99
75) Fluoranthene	18.862	202	449244	20.610	ng	98
77) Benzidine	19.074	184	103924	13.388	ng	96
78) Pyrene	19.227	202	464405	18.471	ng	99
80) Butylbenzylphthalate	20.162	149	132567	12.409	ng	98
81) Benzo(a)anthracene	20.992	228	428024	18.997	ng	99
82) 3,3'-Dichlorobenzidine	20.939	252	120514	19.308	ng	97
83) Chrysene	21.045	228	450797	20.539	ng	99
84) Bis(2-ethylhexyl)phtha...	20.950	149	208257	13.495	ng #	93
85) Di-n-octyl phthalate	21.797	149	341367	14.003	ng	98
87) Indeno(1,2,3-cd)pyrene	25.144	276	400391	19.181	ng	96
88) Benzo(b)fluoranthene	22.486	252	401506	18.085	ng	98
89) Benzo(k)fluoranthene	22.527	252	425028	19.997	ng	99
90) Benzo(a)pyrene	23.015	252	382351	19.629	ng	99
91) Dibenzo(a,h)anthracene	25.156	278	348669	19.259	ng	99
92) Benzo(g,h,i)perylene	25.762	276	343631	20.152	ng	97

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

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