

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090221\
 Data File : BM032028.D
 Acq On : 02 Sep 2021 18:46
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 9/3/2021 3:31:18 PM

Quant Time: Sep 02 19:15:56 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 02 17:39:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.492	152	80578	20.000	ng	0.00
21) Naphthalene-d8	10.251	136	284489	20.000	ng	0.00
39) Acenaphthene-d10	14.139	164	183557	20.000	ng	0.00
64) Phenanthrene-d10	16.892	188	375035	20.000	ng	0.00
76) Chrysene-d12	21.103	240	385787	20.000	ng	0.00
86) Perylene-d12	23.233	264	402919	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.134	112	687794	136.335	ng	0.00
7) Phenol-d6	6.698	99	919421	127.101	ng	0.00
23) Nitrobenzene-d5	8.651	82	912982	145.165	ng	0.00
42) 2,4,6-Tribromophenol	15.645	330	406518	205.161	ng	0.00
45) 2-Fluorobiphenyl	12.751	172	2235442	164.702	ng	0.00
79) Terphenyl-d14	19.545	244	3295699	137.333	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.152	88	140110	69.923	ng	98
3) Pyridine	3.528	79	368824m	70.099	ng	
4) n-Nitrosodimethylamine	3.451	42	163756	57.869	ng	# 71
6) Aniline	6.851	93	472812	61.161	ng	94
8) 2-Chlorophenol	7.069	128	382784	65.515	ng	92
9) Benzaldehyde	6.663	77	212443m	47.883	ng	
10) Phenol	6.728	94	450753	62.809	ng	90
11) bis(2-Chloroethyl)ether	6.951	93	333640	60.913	ng	96
12) 1,3-Dichlorobenzene	7.381	146	437434	70.581	ng	# 92
13) 1,4-Dichlorobenzene	7.528	146	444456	69.885	ng	96
14) 1,2-Dichlorobenzene	7.839	146	430283	69.932	ng	96
15) Benzyl Alcohol	7.751	79	339062	62.165	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.034	45	432706	57.538	ng	97
17) 2-Methylphenol	7.945	107	301881	62.139	ng	96
18) Hexachloroethane	8.545	117	162601	65.510	ng	# 84
19) n-Nitroso-di-n-propyla...	8.310	70	293941	57.345	ng	# 89
20) 3+4-Methylphenols	8.281	107	411379	61.301	ng	95
22) Acetophenone	8.322	105	568782	73.591	ng	# 91
24) Nitrobenzene	8.692	77	455467	71.238	ng	93
25) Isophorone	9.216	82	760580	68.722	ng	# 95
26) 2-Nitrophenol	9.386	139	216817	81.948	ng	# 79
27) 2,4-Dimethylphenol	9.457	122	299407	74.100	ng	97
28) bis(2-Chloroethoxy)met...	9.698	93	409375	70.341	ng	99
29) 2,4-Dichlorophenol	9.916	162	392788	86.619	ng	96
30) 1,2,4-Trichlorobenzene	10.122	180	446455	90.210	ng	97
31) Naphthalene	10.304	128	1179572	74.881	ng	100
32) Benzoic acid	9.669	122	249132	76.805	ng	87
33) 4-Chloroaniline	10.433	127	471930	73.661	ng	# 94
34) Hexachlorobutadiene	10.575	225	281090	93.728	ng	97
35) Caprolactam	11.263	113	103913	73.091	ng	# 78
36) 4-Chloro-3-methylphenol	11.569	107	375889	70.843	ng	90
37) 2-Methylnaphthalene	11.922	142	884990	77.505	ng	95
38) 1-Methylnaphthalene	12.145	142	825945	75.790	ng	97
40) 1,2,4,5-Tetrachloroben...	12.298	216	474217	89.487	ng	# 95
41) Hexachlorocyclopentadiene	12.269	237	248020	101.017	ng	98
43) 2,4,6-Trichlorophenol	12.557	196	325823	86.452	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.627	196	348510	85.124	ng	# 90
46) 1,1'-Biphenyl	12.963	154	1120298	77.936	ng	98
47) 2-Chloronaphthalene	12.998	162	906083	79.974	ng	94
48) 2-Nitroaniline	13.227	65	247848	68.446	ng	# 84
49) Acenaphthylene	13.857	152	1363660	76.183	ng	100
50) Dimethylphthalate	13.621	163	1059738	74.713	ng	99
51) 2,6-Dinitrotoluene	13.739	165	254170	79.876	ng	# 66
52) Acenaphthene	14.204	154	844902	76.991	ng	97
53) 3-Nitroaniline	14.074	138	236115	75.450	ng	85
54) 2,4-Dinitrophenol	14.292	184	160342	90.283	ng	# 63
55) Dibenzofuran	14.545	168	1392095	78.025	ng	92
56) 4-Nitrophenol	14.398	139	196973	78.644	ng	# 75
57) 2,4-Dinitrotoluene	14.539	165	351104	81.231	ng	# 73
58) Fluorene	15.198	166	1130782	78.322	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.774	232	303043	87.423	ng	# 82
60) Diethylphthalate	14.992	149	1061100	71.461	ng	96
61) 4-Chlorophenyl-phenyle...	15.198	204	568554	81.694	ng	90
62) 4-Nitroaniline	15.251	138	229919	75.082	ng	# 80
63) Azobenzene	15.492	77	1041174	63.573	ng	89
65) 4,6-Dinitro-2-methylph...	15.304	198	220284	88.055	ng	# 74
66) n-Nitrosodiphenylamine	15.421	169	965766	76.473	ng	96
67) 4-Bromophenyl-phenylether	16.092	248	344140	86.225	ng	90
68) Hexachlorobenzene	16.198	284	384253	90.087	ng	# 81
69) Atrazine	16.392	200	290733	73.715	ng	94
70) Pentachlorophenol	16.551	266	245425	92.030	ng	98
71) Phenanthrene	16.939	178	1714651	76.973	ng	99
72) Anthracene	17.027	178	1694109	78.021	ng	100
73) Carbazole	17.309	167	1606886	77.012	ng	96
74) Di-n-butylphthalate	17.880	149	1780454	71.066	ng	# 98
75) Fluoranthene	18.962	202	1979961	81.671	ng	98
77) Benzidine	19.168	184	539085	56.938	ng	99
78) Pyrene	19.327	202	2063466	67.290	ng	99
80) Butylbenzylphthalate	20.250	149	804413	61.734	ng	90
81) Benzo(a)anthracene	21.086	228	2017433	73.412	ng	99
82) 3,3'-Dichlorobenzidine	21.033	252	589917	77.489	ng	# 98
83) Chrysene	21.139	228	1977161	73.859	ng	100
84) Bis(2-ethylhexyl)phtha...	21.033	149	1178862	62.630	ng	# 94
85) Di-n-octyl phthalate	21.891	149	1994588	67.081	ng	# 91
87) Indeno(1,2,3-cd)pyrene	25.350	276	2537615	90.537	ng	97
88) Benzo(b)fluoranthene	22.603	252	2192701	73.554	ng	99
89) Benzo(k)fluoranthene	22.650	252	2038096	71.414	ng	98
90) Benzo(a)pyrene	23.144	252	1979325	75.677	ng	99
91) Dibenzo(a,h)anthracene	25.356	278	2193946	90.253	ng	98
92) Benzo(g,h,i)perylene	25.985	276	2105234	91.947	ng	99

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

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Abundance

TIC: BM032028.D\data.ms

Time-->