

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040121\
 Data File : BM029302.D
 Acq On : 01 Apr 2021 12:25
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 4/2/2021 12:08:48 PM

Quant Time: Apr 01 14:17:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 14:16:29 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	99448	20.00	ng	0.00
21) Naphthalene-d8	10.486	136	384722	20.00	ng	0.00
39) Acenaphthene-d10	14.339	164	230517	20.00	ng	0.00
64) Phenanthrene-d10	17.074	188	443218	20.00	ng	0.00
76) Chrysene-d12	21.250	240	418758	20.00	ng	0.00
86) Perylene-d12	23.462	264	442825	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.298	112	125894	21.82	ng	0.00
7) Phenol-d6	6.875	99	174860	21.12	ng	0.00
23) Nitrobenzene-d5	8.851	82	166341	20.99	ng	0.00
42) 2,4,6-Tribromophenol	15.827	330	43201	20.24	ng	0.00
45) 2-Fluorobiphenyl	12.957	172	329354	20.27	ng	0.00
79) Terphenyl-d14	19.703	244	433360	20.79	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.234	88	28519	10.88	ng	99
3) Pyridine	3.640	79	69345	10.97	ng	95
4) n-Nitrosodimethylamine	3.551	42	31011	10.93	ng	# 92
6) Aniline	7.034	93	95469	10.64	ng	99
8) 2-Chlorophenol	7.269	128	68701	10.55	ng	97
9) Benzaldehyde	6.851	77	59273m	11.40	ng	
10) Phenol	6.904	94	86172	10.54	ng	96
11) bis(2-Chloroethyl)ether	7.134	93	66251	11.32	ng	98
12) 1,3-Dichlorobenzene	7.592	146	75483	10.36	ng	96
13) 1,4-Dichlorobenzene	7.739	146	78091	10.57	ng	95
14) 1,2-Dichlorobenzene	8.051	146	73991	10.43	ng	99
15) Benzyl Alcohol	7.945	79	61792	10.32	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.233	45	99465	11.08	ng	99
17) 2-Methylphenol	8.145	107	56034	10.66	ng	96
18) Hexachloroethane	8.775	117	30604	11.41	ng	95
19) n-Nitroso-di-n-propyla...	8.504	70	57985	10.86	ng	99
20) 3+4-Methylphenols	8.469	107	74898	10.65	ng	96
22) Acetophenone	8.522	105	100810	10.47	ng	99
24) Nitrobenzene	8.898	77	86672	10.48	ng	97
25) Isophorone	9.422	82	149365	10.97	ng	99
26) 2-Nitrophenol	9.604	139	32698	11.86	ng	90
27) 2,4-Dimethylphenol	9.669	122	53269	10.08	ng	96
28) bis(2-Chloroethoxy)met...	9.904	93	77780	10.81	ng	99
29) 2,4-Dichlorophenol	10.133	162	58909	10.25	ng	96
30) 1,2,4-Trichlorobenzene	10.351	180	65386	10.02	ng	99
31) Naphthalene	10.533	128	206076	10.16	ng	99
32) Benzoic acid	9.751	122	32377	10.54	ng	95
33) 4-Chloroaniline	10.645	127	81493	10.17	ng	97
34) Hexachlorobutadiene	10.822	225	39163	10.33	ng	92
35) Caprolactam	11.416	113	17766	10.18	ng	93
36) 4-Chloro-3-methylphenol	11.769	107	62411	10.17	ng	94
37) 2-Methylnaphthalene	12.151	142	143389	10.11	ng	98
38) 1-Methylnaphthalene	12.369	142	136293	10.10	ng	99
40) 1,2,4,5-Tetrachloroben...	12.521	216	63520	9.83	ng	98
41) Hexachlorocyclopentadiene	12.504	237	36047	10.17	ng	99
43) 2,4,6-Trichlorophenol	12.763	196	44714	10.63	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.833	196	48957	10.47	ng	97
46) 1,1'-Biphenyl	13.168	154	180755	10.31	ng	98
47) 2-Chloronaphthalene	13.210	162	137009	9.93	ng	99
48) 2-Nitroaniline	13.416	65	43896	10.85	ng	100
49) Acenaphthylene	14.057	152	211813	10.09	ng	99
50) Dimethylphthalate	13.798	163	166004	10.31	ng	100
51) 2,6-Dinitrotoluene	13.910	165	36652	10.52	ng	89
52) Acenaphthene	14.404	154	132597	9.94	ng	98
53) 3-Nitroaniline	14.239	138	36360	10.31	ng	94
54) 2,4-Dinitrophenol	14.445	184	14952	8.59	ng	91
55) Dibenzofuran	14.733	168	211112	10.06	ng	98
56) 4-Nitrophenol	14.551	139	28608	8.88	ng	99
57) 2,4-Dinitrotoluene	14.704	165	46170	10.08	ng	99
58) Fluorene	15.386	166	168890	9.93	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.962	232	36254	9.46	ng	98
60) Diethylphthalate	15.162	149	169249	10.67	ng	99
61) 4-Chlorophenyl-phenyle...	15.386	204	77211	9.84	ng	97
62) 4-Nitroaniline	15.398	138	37089	9.99	ng	99
63) Azobenzene	15.674	77	191347	10.55	ng	98
65) 4,6-Dinitro-2-methylph...	15.462	198	22503	9.11	ng	92
66) n-Nitrosodiphenylamine	15.598	169	146969	10.36	ng	100
67) 4-Bromophenyl-phenylether	16.274	248	43452	10.69	ng	97
68) Hexachlorobenzene	16.386	284	47704	9.98	ng	98
69) Atrazine	16.551	200	47976	10.71	ng	98
70) Pentachlorophenol	16.733	266	16197	5.50	ng	96
71) Phenanthrene	17.121	178	257173	10.01	ng	99
72) Anthracene	17.209	178	250048	10.01	ng	99
73) Carbazole	17.480	167	245170	9.98	ng	98
74) Di-n-butylphthalate	18.051	149	270345	11.41	ng	99
75) Fluoranthene	19.133	202	282354	9.94	ng	99
77) Benzidine	19.321	184	121458	10.04	ng	99
78) Pyrene	19.492	202	294471	10.39	ng	100
80) Butylbenzylphthalate	20.397	149	129920	13.56	ng	95
81) Benzo(a)anthracene	21.233	228	277438	10.15	ng	99
82) 3,3'-Dichlorobenzidine	21.174	252	91593	11.21	ng	95
83) Chrysene	21.286	228	274416	10.04	ng	98
84) Bis(2-ethylhexyl)phtha...	21.174	149	196100	14.24	ng	99
85) Di-n-octyl phthalate	22.044	149	340939	15.28	ng	97
87) Indeno(1,2,3-cd)pyrene	25.691	276	324776	10.03	ng	98
88) Benzo(b)fluoranthene	22.797	252	294283	10.22	ng	99
89) Benzo(k)fluoranthene	22.844	252	271630	9.70	ng	99
90) Benzo(a)pyrene	23.368	252	264547	10.14	ng	99
91) Dibenzo(a,h)anthracene	25.703	278	279408	10.08	ng	97
92) Benzo(g,h,i)perylene	26.373	276	274712	10.11	ng	98

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

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