

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041724\
 Data File : BG060919.D
 Acq On : 17 Apr 2024 13:51
 Operator : MA/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC020

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 04/22/2024

Supervised By :mohammad ahmed 04/22/2024

Quant Time: Apr 18 01:52:59 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG041724.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Apr 18 01:47:49 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	40524	20.000	ng	0.00
21) Naphthalene-d8	10.738	136	164435	20.000	ng	0.00
39) Acenaphthene-d10	14.568	164	142039	20.000	ng	0.00
64) Phenanthrene-d10	17.318	188	371535	20.000	ng	0.00
76) Chrysene-d12	21.578	240	358960	20.000	ng	0.00
86) Perylene-d12	24.698	264	422652	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	96976	41.567	ng	0.00
7) Phenol-d6	7.107	99	141233	42.022	ng	0.00
23) Nitrobenzene-d5	9.092	82	152978	42.235	ng	0.00
42) 2,4,6-Tribromophenol	16.061	330	98984	42.183	ng	0.00
45) 2-Fluorobiphenyl	13.194	172	402873	41.622	ng	0.00
79) Terphenyl-d14	19.939	244	916409	44.555	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.452	88	21007	22.274	ng	95
3) Pyridine	3.840	79	62320m	20.795	ng	
4) n-Nitrosodimethylamine	3.758	42	39599	20.681	ng	88
6) Aniline	7.265	93	88888	21.707	ng	97
8) 2-Chlorophenol	7.512	128	56138	21.151	ng	98
9) Benzaldehyde	7.083	77	42503	23.575	ng	95
10) Phenol	7.130	94	69310	20.577	ng	93
11) bis(2-Chloroethyl)ether	7.365	93	54963	21.037	ng	97
12) 1,3-Dichlorobenzene	7.835	146	59857	21.226	ng	96
13) 1,4-Dichlorobenzene	7.976	146	60497	20.955	ng	96
14) 1,2-Dichlorobenzene	8.293	146	59455	21.178	ng	93
15) Benzyl Alcohol	8.170	79	60465	20.914	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.476	45	130622m	20.687	ng	
17) 2-Methylphenol	8.382	107	55564	20.878	ng	98
18) Hexachloroethane	9.022	117	21200	20.680	ng	96
19) n-Nitroso-di-n-propyla...	8.740	70	53478	20.890	ng	99
20) 3+4-Methylphenols	8.705	107	78501	21.261	ng	98
22) Acetophenone	8.758	105	98763	21.091	ng	# 98
24) Nitrobenzene	9.134	77	76097	21.407	ng	92
25) Isophorone	9.662	82	143454	21.462	ng	98
26) 2-Nitrophenol	9.845	139	31097	20.823	ng	91
27) 2,4-Dimethylphenol	9.909	122	53498	20.332	ng	95
28) bis(2-Chloroethoxy)met...	10.144	93	78370	20.851	ng	98
29) 2,4-Dichlorophenol	10.385	162	60622	21.235	ng	98
30) 1,2,4-Trichlorobenzene	10.602	180	64203	20.981	ng	99
31) Naphthalene	10.790	128	187058	21.034	ng	99
32) Benzoic acid	10.009	122	46183m	21.654	ng	
33) 4-Chloroaniline	10.890	127	87754	21.055	ng	98
34) Hexachlorobutadiene	11.078	225	45014	20.692	ng	94
35) Caprolactam	11.637	113	22746	21.595	ng	# 86
36) 4-Chloro-3-methylphenol	12.013	107	75669	21.931	ng	97
37) 2-Methylnaphthalene	12.394	142	143576	21.341	ng	99
38) 1-Methylnaphthalene	12.612	142	137089	21.585	ng	99
40) 1,2,4,5-Tetrachloroben...	12.765	216	89236	20.359	ng	99
41) Hexachlorocyclopentadiene	12.747	237	43811	19.548	ng	98
43) 2,4,6-Trichlorophenol	13.000	196	61848	20.292	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.070	196	76148	20.717	ng	94
46) 1,1'-Biphenyl	13.405	154	199910	20.699	ng	97
47) 2-Chloronaphthalene	13.446	162	154175	20.858	ng	98
48) 2-Nitroaniline	13.640	65	63259	20.855	ng	99
49) Acenaphthylene	14.292	152	258124	20.651	ng	99
50) Dimethylphthalate	14.028	163	237498	21.071	ng	99
51) 2,6-Dinitrotoluene	14.139	165	48722	21.209	ng	95
52) Acenaphthene	14.633	154	158992m	20.974	ng	
53) 3-Nitroaniline	14.469	138	54255	21.514	ng	# 98
54) 2,4-Dinitrophenol	14.674	184	22335	19.620	ng	95
55) Dibenzofuran	14.974	168	270263	21.105	ng	98
56) 4-Nitrophenol	14.774	139	39361	20.386	ng	92
57) 2,4-Dinitrotoluene	14.927	165	70333	21.425	ng	99
58) Fluorene	15.620	166	228108	21.475	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.197	232	71325	21.123	ng	98
60) Diethylphthalate	15.397	149	239918	21.251	ng	99
61) 4-Chlorophenyl-phenyle...	15.614	204	126230	21.439	ng	100
62) 4-Nitroaniline	15.626	138	57632	21.255	ng	96
63) Azobenzene	15.908	77	225527	21.196	ng	98
65) 4,6-Dinitro-2-methylph...	15.691	198	39724	19.164	ng	100
66) n-Nitrosodiphenylamine	15.826	169	213677	20.951	ng	99
67) 4-Bromophenyl-phenylether	16.513	248	86548	20.794	ng	97
68) Hexachlorobenzene	16.625	284	94021	20.760	ng	99
69) Atrazine	16.778	200	79052	21.567	ng	98
70) Pentachlorophenol	16.966	266	62896	20.507	ng	98
71) Phenanthrene	17.359	178	400770	21.311	ng	98
72) Anthracene	17.453	178	413710	21.417	ng	99
73) Carbazole	17.718	167	356614	21.189	ng	99
74) Di-n-butylphthalate	18.293	149	434784	21.458	ng	99
75) Fluoranthene	19.375	202	516864	21.593	ng	98
77) Benzidine	19.551	184	144818	18.290	ng	98
78) Pyrene	19.733	202	531640	21.590	ng	98
80) Butylbenzylphthalate	20.632	149	190844	21.101	ng	98
81) Benzo(a)anthracene	21.560	228	536920	21.182	ng	99
82) 3,3'-Dichlorobenzidine	21.478	252	191577	21.224	ng	99
83) Chrysene	21.625	228	509883	21.068	ng	99
84) Bis(2-ethylhexyl)phtha...	21.490	149	286281	21.406	ng	100
85) Di-n-octyl phthalate	22.682	149	487070	21.107	ng	100
87) Indeno(1,2,3-cd)pyrene	28.241	276	616204	21.083	ng	# 93
88) Benzo(b)fluoranthene	23.716	252	499726	20.897	ng	100
89) Benzo(k)fluoranthene	23.781	252	511895	20.925	ng	98
90) Benzo(a)pyrene	24.551	252	492764	20.900	ng	99
91) Dibenzo(a,h)anthracene	28.305	278	510918	21.237	ng	97
92) Benzo(g,h,i)perylene	29.339	276	479420	20.383	ng	98

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

