

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120723\
 Data File : BG059989.D
 Acq On : 7 Dec 2023 16:42
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

ICVBG120723

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 12/08/2023

Supervised By :mohammad ahmed 12/08/2023

Quant Time: Dec 08 02:37:23 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG120723.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Dec 08 02:35:41 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.959	152	129695	20.000	ng	0.00
21) Naphthalene-d8	10.768	136	435977	20.000	ng	# 0.00
39) Acenaphthene-d10	14.598	164	421338	20.000	ng	0.00
64) Phenanthrene-d10	17.342	188	1411934	20.000	ng	0.00
76) Chrysene-d12	21.620	240	1923884	20.000	ng	0.00
86) Perylene-d12	24.810	264	2279617	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.527	112	530072	75.584	ng	0.00
7) Phenol-d6	7.113	99	725919	73.623	ng	0.00
23) Nitrobenzene-d5	9.122	82	797333	80.763	ng	0.00
42) 2,4,6-Tribromophenol	16.085	330	1002007	81.818	ng	0.00
45) 2-Fluorobiphenyl	13.223	172	2850328	81.291	ng	0.00
79) Terphenyl-d14	19.969	244	6208150	82.296	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.441	88	104097	32.819	ng	98
3) Pyridine	3.829	79	333544	37.905	ng	95
4) n-Nitrosodimethylamine	3.746	42	147402	38.640	ng	# 63
6) Aniline	7.283	93	446350	42.466	ng	97
8) 2-Chlorophenol	7.518	128	241668	35.899	ng	94
9) Benzaldehyde	7.095	77	228743	38.036	ng	94
10) Phenol	7.142	94	354846	36.834	ng	97
11) bis(2-Chloroethyl)ether	7.383	93	217200	33.704	ng	96
12) 1,3-Dichlorobenzene	7.847	146	303910m	31.933	ng	
13) 1,4-Dichlorobenzene	7.994	146	317307	33.886	ng	99
14) 1,2-Dichlorobenzene	8.312	146	314299	35.191	ng	95
15) Benzyl Alcohol	8.194	79	338630	38.485	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.482	45	120079	34.598	ng	91
17) 2-Methylphenol	8.388	107	257517	39.548	ng	95
18) Hexachloroethane	9.046	117	105888	34.848	ng	90
19) n-Nitroso-di-n-propyla...	8.764	70	244544	36.807	ng	98
20) 3+4-Methylphenols	8.723	107	350411	37.993	ng	96
22) Acetophenone	8.782	105	490151	38.433	ng	# 95
24) Nitrobenzene	9.164	77	394780	38.710	ng	95
25) Isophorone	9.686	82	680580	38.325	ng	99
26) 2-Nitrophenol	9.874	139	147340	37.155	ng	91
27) 2,4-Dimethylphenol	9.927	122	263659	49.276	ng	94
28) bis(2-Chloroethoxy)met...	10.168	93	325513	38.699	ng	98
29) 2,4-Dichlorophenol	10.403	162	391620	38.847	ng	97
30) 1,2,4-Trichlorobenzene	10.627	180	524486	37.466	ng	98
31) Naphthalene	10.815	128	863613	38.782	ng	97
32) Benzoic acid	10.057	122	215881	39.902	ng	85
33) 4-Chloroaniline	10.920	127	364314	43.474	ng	95
34) Hexachlorobutadiene	11.102	225	562081	36.876	ng	92
35) Caprolactam	11.696	113	87591	38.173	ng	90
36) 4-Chloro-3-methylphenol	12.037	107	343068	38.798	ng	99
37) 2-Methylnaphthalene	12.424	142	752891	39.031	ng	99
38) 1-Methylnaphthalene	12.642	142	720828	39.341	ng	94
40) 1,2,4,5-Tetrachloroben...	12.789	216	1088297	40.388	ng	98
41) Hexachlorocyclopentadiene	12.771	237	595052	47.261	ng	96
43) 2,4,6-Trichlorophenol	13.024	196	548572	39.282	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.094	196	653799	41.516	ng	99
46) 1,1'-Biphenyl	13.429	154	1190600	38.848	ng	96
47) 2-Chloronaphthalene	13.470	162	1000273	37.908	ng	98
48) 2-Nitroaniline	13.670	65	219876	42.683	ng	86
49) Acenaphthylene	14.316	152	1441293	40.511	ng	100
50) Dimethylphthalate	14.052	163	1377546	41.018	ng	99
51) 2,6-Dinitrotoluene	14.169	165	288278	38.996	ng	93
52) Acenaphthene	14.663	154	1019377	39.521	ng	99
53) 3-Nitroaniline	14.498	138	199994	39.557	ng	# 94
54) 2,4-Dinitrophenol	14.698	184	252531	41.097	ng	97
55) Dibenzofuran	14.998	168	1692163	40.380	ng	100
56) 4-Nitrophenol	14.792	139	166181	38.795	ng	99
57) 2,4-Dinitrotoluene	14.951	165	422308	39.892	ng	95
58) Fluorene	15.644	166	1440674	39.458	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.215	232	713012	40.358	ng	98
60) Diethylphthalate	15.421	149	1219580	40.387	ng	99
61) 4-Chlorophenyl-phenyle...	15.638	204	1231491	40.175	ng	96
62) 4-Nitroaniline	15.662	138	227590	40.633	ng	96
63) Azobenzene	15.932	77	1041266	39.469	ng	96
65) 4,6-Dinitro-2-methylph...	15.715	198	397990	39.543	ng	99
66) n-Nitrosodiphenylamine	15.850	169	1369324	40.819	ng	99
67) 4-Bromophenyl-phenylether	16.537	248	909442	41.055	ng	96
68) Hexachlorobenzene	16.649	284	958235	37.190	ng	97
69) Atrazine	16.802	200	461188	31.499	ng	97
70) Pentachlorophenol	16.990	266	709170	39.320	ng	98
71) Phenanthrene	17.389	178	2661205	41.170	ng	99
72) Anthracene	17.477	178	2742397	41.979	ng	99
73) Carbazole	17.748	167	2061091	38.689	ng	# 93
74) Di-n-butylphthalate	18.312	149	1902545	39.973	ng	99
75) Fluoranthene	19.404	202	3874684	38.285	ng	97
77) Benzidine	19.581	184	649226	28.763	ng	99
78) Pyrene	19.763	202	3937730	38.212	ng	98
80) Butylbenzylphthalate	20.662	149	719325	39.597	ng	95
81) Benzo(a)anthracene	21.596	228	4762250	38.784	ng	98
82) 3,3'-Dichlorobenzidine	21.514	252	1882086	42.988	ng	97
83) Chrysene	21.667	228	4507478	38.884	ng	100
84) Bis(2-ethylhexyl)phtha...	21.520	149	1105299	41.335	ng	97
85) Di-n-octyl phthalate	22.724	149	1866203	40.381	ng	100
87) Indeno(1,2,3-cd)pyrene	28.459	276	6468852	38.950	ng	100
88) Benzo(b)fluoranthene	23.799	252	5023177m	37.868	ng	
89) Benzo(k)fluoranthene	23.864	252	5175898	40.278	ng	99
90) Benzo(a)pyrene	24.663	252	5002036	41.361	ng	# 98
91) Dibenzo(a,h)anthracene	28.523	278	5558962	39.739	ng	98
92) Benzo(g,h,i)perylene	29.593	276	5254717	38.546	ng	# 97

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

