

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
 Data File : BG060240.D
 Acq On : 4 Jan 2024 11:36
 Operator : MA/JU
 Sample : PB158168BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB158168BS

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 01/05/2024

Supervised By :mohammad ahmed 01/05/2024

Quant Time: Jan 04 12:34:28 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Dec 30 03:14:04 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.935	152	170429	20.000	ng	0.00
21) Naphthalene-d8	10.743	136	767179	20.000	ng	0.00
39) Acenaphthene-d10	14.574	164	610194	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	1559266	20.000	ng	0.00
76) Chrysene-d12	21.595	240	1417944	20.000	ng	0.00
86) Perylene-d12	24.768	264	1508731	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.514	112	1577881	139.739	ng	0.00
7) Phenol-d6	7.106	99	2370295	141.380	ng	0.00
23) Nitrobenzene-d5	9.104	82	1330834	80.126	ng	0.00
42) 2,4,6-Tribromophenol	16.066	330	1110735	135.985	ng	0.00
45) 2-Fluorobiphenyl	13.199	172	3332593	78.494	ng	0.00
79) Terphenyl-d14	19.944	244	4978803	67.504	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.411	88	164396	32.724	ng	100
3) Pyridine	3.804	79	406597	28.538	ng	96
4) n-Nitrosodimethylamine	3.722	42	204803	38.533	ng	# 93
6) Aniline	7.265	93	524916	31.249	ng	95
8) 2-Chlorophenol	7.506	128	519873	47.906	ng	98
9) Benzaldehyde	7.077	77	387410	38.434	ng	94
10) Phenol	7.130	94	824633	48.776	ng	97
11) bis(2-Chloroethyl)ether	7.359	93	552884	40.413	ng	94
12) 1,3-Dichlorobenzene	7.829	146	529197	39.504	ng	99
13) 1,4-Dichlorobenzene	7.976	146	538490	39.761	ng	93
14) 1,2-Dichlorobenzene	8.293	146	530894	40.097	ng	99
15) Benzyl Alcohol	8.176	79	526348m	50.579	ng	
16) 2,2'-oxybis(1-Chloropr...	8.464	45	735355m	41.881	ng	
17) 2-Methylphenol	8.381	107	533216	50.368	ng	99
18) Hexachloroethane	9.022	117	178689	41.893	ng	95
19) n-Nitroso-di-n-propyla...	8.740	70	496506	43.350	ng	97
20) 3+4-Methylphenols	8.710	107	733876	48.912	ng	96
22) Acetophenone	8.757	105	894256	39.342	ng	# 97
24) Nitrobenzene	9.145	77	684622	40.287	ng	99
25) Isophorone	9.662	82	1363565	39.239	ng	98
26) 2-Nitrophenol	9.856	139	282435	47.629	ng	98
27) 2,4-Dimethylphenol	9.909	122	477518	56.849	ng	97
28) bis(2-Chloroethoxy)met...	10.150	93	829456	40.754	ng	97
29) 2,4-Dichlorophenol	10.391	162	624341	47.030	ng	95
30) 1,2,4-Trichlorobenzene	10.602	180	615886	38.117	ng	98
31) Naphthalene	10.796	128	1604849	39.624	ng	98
32) Benzoic acid	10.044	122	373567	49.388	ng	94
33) 4-Chloroaniline	10.902	127	384402	24.302	ng	97
34) Hexachlorobutadiene	11.072	225	341604	35.548	ng	95
35) Caprolactam	11.666	113	223476m	50.607	ng	
36) 4-Chloro-3-methylphenol	12.024	107	659204	48.250	ng	98
37) 2-Methylnaphthalene	12.400	142	1216393	41.579	ng	98
38) 1-Methylnaphthalene	12.623	142	1190891	40.263	ng	99
40) 1,2,4,5-Tetrachloroben...	12.770	216	753864	38.569	ng	99
41) Hexachlorocyclopentadiene	12.747	237	702681	111.224	ng	99
43) 2,4,6-Trichlorophenol	13.005	196	579475	45.115	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.082	196	644326	44.843	ng	98
46) 1,1'-Biphenyl	13.411	154	1864844	41.539	ng	98
47) 2-Chloronaphthalene	13.452	162	1536058	40.160	ng	98
48) 2-Nitroaniline	13.658	65	499271	49.948	ng	98
49) Acenaphthylene	14.298	152	2500061	45.365	ng	98
50) Dimethylphthalate	14.034	163	2038107	43.261	ng	99
51) 2,6-Dinitrotoluene	14.151	165	443509	45.260	ng	96
52) Acenaphthene	14.645	154	1413757	41.110	ng	97
53) 3-Nitroaniline	14.480	138	313295	33.931	ng	97
54) 2,4-Dinitrophenol	14.686	184	548000	94.382	ng	99
55) Dibenzofuran	14.974	168	2456141	42.637	ng	99
56) 4-Nitrophenol	14.791	139	776429	111.247	ng	96
57) 2,4-Dinitrotoluene	14.938	165	632751	47.488	ng	# 99
58) Fluorene	15.626	166	2012931	42.605	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.203	232	583143	46.563	ng	98
60) Diethylphthalate	15.397	149	1971491	43.882	ng	98
61) 4-Chlorophenyl-phenyle...	15.614	204	993844	40.398	ng	97
62) 4-Nitroaniline	15.649	138	454692	46.723	ng	95
63) Azobenzene	15.914	77	2005100	43.048	ng	98
65) 4,6-Dinitro-2-methylph...	15.702	198	375067	48.072	ng	98
66) n-Nitrosodiphenylamine	15.831	169	1810263	43.084	ng	98
67) 4-Bromophenyl-phenylether	16.513	248	657225	39.756	ng	96
68) Hexachlorobenzene	16.631	284	757345	39.073	ng	97
69) Atrazine	16.777	200	540336	34.990	ng	97
70) Pentachlorophenol	16.971	266	971546	91.243	ng	99
71) Phenanthrene	17.371	178	3349005	43.452	ng	100
72) Anthracene	17.459	178	3339439	43.620	ng	99
73) Carbazole	17.729	167	3127907	43.077	ng	98
74) Di-n-butylphthalate	18.287	149	3266443	46.421	ng	98
75) Fluoranthene	19.380	202	3763725	41.632	ng	99
77) Benzidine	19.568	184	1201262	30.610	ng	99
78) Pyrene	19.744	202	3910693	41.029	ng	99
80) Butylbenzylphthalate	20.632	149	1505474	53.566	ng	96
81) Benzo(a)anthracene	21.572	228	3842284	42.619	ng	99
82) 3,3'-Dichlorobenzidine	21.490	252	1323404	40.367	ng	97
83) Chrysene	21.636	228	3678352	41.997	ng	99
84) Bis(2-ethylhexyl)phtha...	21.478	149	2241791	52.598	ng	99
85) Di-n-octyl phthalate	22.676	149	3775866	54.121	ng	100
87) Indeno(1,2,3-cd)pyrene	28.399	276	4761597	44.765	ng	99
88) Benzo(b)fluoranthene	23.763	252	4049508	44.664	ng	99
89) Benzo(k)fluoranthene	23.828	252	3995615	43.785	ng	99
90) Benzo(a)pyrene	24.621	252	3628608	44.237	ng	98
91) Dibenzo(a,h)anthracene	28.458	278	3987332	45.622	ng	99
92) Benzo(g,h,i)perylene	29.527	276	3840518	43.383	ng	99

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

