

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121622\
 Data File : BG055998.D
 Acq On : 16 Dec 2022 17:47
 Operator : CG/JU
 Sample : N6037-19MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RB-25-(7-9)MSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Dec 17 01:22:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 17 01:18:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.376	152	8852	20.000	ng	0.00
21) Naphthalene-d8	11.208	136	32226	20.000	ng	# 0.00
39) Acenaphthene-d10	14.980	164	27730	20.000	ng	0.00
64) Phenanthrene-d10	17.718	188	80173	20.000	ng	# 0.00
76) Chrysene-d12	22.030	240	78147	20.000	ng	0.00
86) Perylene-d12	25.526	264	95512	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.890	112	43300	114.716	ng	0.00
7) Phenol-d6	7.506	99	53401	110.001	ng	0.00
23) Nitrobenzene-d5	9.545	82	33677	73.029	ng	0.00
42) 2,4,6-Tribromophenol	16.460	330	83782	125.300	ng	0.00
45) 2-Fluorobiphenyl	13.611	172	133664	67.602	ng	0.00
79) Terphenyl-d14	20.291	244	320590	72.505	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.716	88	4440	40.420	ng	99
3) Pyridine	4.133	79	13780	38.119	ng	95
4) n-Nitrosodimethylamine	4.039	42	12665	38.974	ng	# 99
6) Aniline	7.682	93	19753	33.985	ng	96
8) 2-Chlorophenol	7.929	128	22084	45.350	ng	94
9) Benzaldehyde	7.494	77	13139	43.095	ng	89
10) Phenol	7.529	94	23261	43.780	ng	98
11) bis(2-Chloroethyl)ether	7.770	93	14550	42.042	ng	93
12) 1,3-Dichlorobenzene	8.264	146	26530	42.116	ng	90
13) 1,4-Dichlorobenzene	8.411	146	26641	41.382	ng	97
14) 1,2-Dichlorobenzene	8.734	146	25881	42.066	ng	95
15) Benzyl Alcohol	8.605	79	20982	41.595	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.899	45	19539	43.379	ng	97
17) 2-Methylphenol	8.799	107	18399	44.013	ng	94
18) Hexachloroethane	9.474	117	8560	43.804	ng	# 87
19) n-Nitroso-di-n-propyla...	9.181	70	13887	39.908	ng	# 88
20) 3+4-Methylphenols	9.128	107	25529	42.794	ng	97
22) Acetophenone	9.198	105	32784	41.038	ng	# 96
24) Nitrobenzene	9.586	77	22424	48.310	ng	97
25) Isophorone	10.115	82	40077	41.445	ng	97
26) 2-Nitrophenol	10.309	139	13702	49.865	ng	97
27) 2,4-Dimethylphenol	10.344	122	23088m	50.031	ng	
28) bis(2-Chloroethoxy)met...	10.585	93	19953	46.688	ng	97
29) 2,4-Dichlorophenol	10.837	162	29356	44.818	ng	96
30) 1,2,4-Trichlorobenzene	11.061	180	35619	41.833	ng	94
31) Naphthalene	11.260	128	69111	40.644	ng	98
32) Benzoic acid	10.473	122	19096	48.244	ng	92
33) 4-Chloroaniline	11.354	127	13597	18.653	ng	96
34) Hexachlorobutadiene	11.537	225	31597	39.472	ng	94
35) Caprolactam	12.124	113	7237m	39.656	ng	
36) 4-Chloro-3-methylphenol	12.441	107	26478	43.111	ng	97
37) 2-Methylnaphthalene	12.835	142	59825	39.823	ng	96
38) 1-Methylnaphthalene	13.052	142	54846	37.476	ng	99
40) 1,2,4,5-Tetrachloroben...	13.193	216	54124	46.261	ng	99
41) Hexachlorocyclopentadiene	13.170	237	72030	108.522	ng	98
43) 2,4,6-Trichlorophenol	13.423	196	34134	49.286	ng	95

12/19/2022
 Supervised By :mohammad
 Ahmed

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.493	196	37823	48.447	ng	99	12/19/2022
46) 1,1'-Biphenyl	13.822	154	87202	46.394	ng	99	Supervised By :mohammad
47) 2-Chloronaphthalene	13.869	162	69058	45.217	ng	97	ahmed
48) 2-Nitroaniline	14.057	65	16531	54.502	ng	90	
49) Acenaphthylene	14.709	152	111527	45.749	ng	100	
50) Dimethylphthalate	14.421	163	98970	41.282	ng	99	
51) 2,6-Dinitrotoluene	14.545	165	20408	55.008	ng	92	12/19/2022
52) Acenaphthene	15.044	154	85024	51.952	ng	95	
53) 3-Nitroaniline	14.868	138	11154	29.558	ng	# 86	
54) 2,4-Dinitrophenol	15.074	184	29330	106.343	ng	# 86	
55) Dibenzofuran	15.373	168	120701	43.858	ng	96	
56) 4-Nitrophenol	15.156	139	31354	103.684	ng	96	
57) 2,4-Dinitrotoluene	15.326	165	29690	53.017	ng	93	
58) Fluorene	16.020	166	97051	43.140	ng	97	
59) 2,3,4,6-Tetrachlorophenol	15.591	232	42784	47.970	ng	99	
60) Diethylphthalate	15.773	149	95375	40.702	ng	98	
61) 4-Chlorophenyl-phenyle...	16.002	204	66861	43.627	ng	98	
62) 4-Nitroaniline	16.031	138	19186	46.686	ng	93	
63) Azobenzene	16.296	77	60967	42.747	ng	99	
65) 4,6-Dinitro-2-methylph...	16.084	198	20063	52.385	ng	94	
66) n-Nitrosodiphenylamine	16.213	169	90072	42.872	ng	98	
67) 4-Bromophenyl-phenylether	16.901	248	50811	43.251	ng	91	
68) Hexachlorobenzene	17.024	284	61632	42.592	ng	97	
69) Atrazine	17.154	200	44072	46.328	ng	100	
70) Pentachlorophenol	17.359	266	79040	95.544	ng	96	
71) Phenanthrene	17.765	178	179349	42.687	ng	98	
72) Anthracene	17.853	178	178525	42.949	ng	100	
73) Carbazole	18.111	167	142508	41.570	ng	98	
74) Di-n-butylphthalate	18.646	149	158354	39.720	ng	100	
75) Fluoranthene	19.750	202	231419	39.335	ng	99	
77) Benzidine	19.915	184	101343	73.844	ng	98	
78) Pyrene	20.109	202	239415	50.630	ng	99	
80) Butylbenzylphthalate	20.978	149	61704	46.050	ng	99	
81) Benzo(a)anthracene	22.007	228	248606	45.209	ng	100	
82) 3,3'-Dichlorobenzidine	21.907	252	53115	26.846	ng	94	
83) Chrysene	22.077	228	237148	46.014	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.883	149	87507	43.957	ng	98	
85) Di-n-octyl phthalate	23.194	149	149356	44.229	ng	99	
87) Indeno(1,2,3-cd)pyrene	29.568	276	323702	42.146	ng	# 98	
88) Benzo(b)fluoranthene	24.410	252	281389	45.522	ng	100	
89) Benzo(k)fluoranthene	24.486	252	260850	42.296	ng	100	
90) Benzo(a)pyrene	25.367	252	245673	47.018	ng	98	
91) Dibenzo(a,h)anthracene	29.645	278	267276	41.577	ng	99	
92) Benzo(g,h,i)perylene	30.843	276	265774	42.760	ng	96	

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

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