

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041624\
 Data File : BG060909.D
 Acq On : 16 Apr 2024 11:36
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 04/17/2024

Supervised By :mohammad ahmed 04/18/2024

Quant Time: Apr 16 19:49:08 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Mar 30 03:12:26 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	87907	20.000	ng	-0.01
21) Naphthalene-d8	10.749	136	346268	20.000	ng	0.00
39) Acenaphthene-d10	14.580	164	266577	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	624499	20.000	ng	0.00
76) Chrysene-d12	21.595	240	575816	20.000	ng	0.00
86) Perylene-d12	24.727	264	674847	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.532	112	375800	71.216	ng	0.00
7) Phenol-d6	7.118	99	530986	67.789	ng	0.00
23) Nitrobenzene-d5	9.110	82	571081	94.109	ng	0.00
42) 2,4,6-Tribromophenol	16.072	330	324045	107.084	ng	0.00
45) 2-Fluorobiphenyl	13.205	172	1382566	76.117	ng	-0.01
79) Terphenyl-d14	19.950	244	2377703	72.710	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.464	88	68245	31.544	ng	98
3) Pyridine	3.845	79	219205	33.590	ng	87
4) n-Nitrosodimethylamine	3.763	42	146283	37.500	ng	# 88
6) Aniline	7.277	93	323349	32.690	ng	97
8) 2-Chlorophenol	7.518	128	212870	35.583	ng	99
9) Benzaldehyde	7.089	77	146472m	33.911	ng	
10) Phenol	7.147	94	265273	33.184	ng	94
11) bis(2-Chloroethyl)ether	7.377	93	208587	32.319	ng	98
12) 1,3-Dichlorobenzene	7.841	146	226137	36.349	ng	97
13) 1,4-Dichlorobenzene	7.982	146	232126	36.572	ng	99
14) 1,2-Dichlorobenzene	8.305	146	227733	36.613	ng	98
15) Benzyl Alcohol	8.187	79	230421	36.316	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.475	45	497862m	34.181	ng	
17) 2-Methylphenol	8.387	107	213336	35.351	ng	94
18) Hexachloroethane	9.028	117	85799	38.062	ng	89
19) n-Nitroso-di-n-propyla...	8.757	70	194340	32.534	ng	94
20) 3+4-Methylphenols	8.722	107	289859	35.076	ng	92
22) Acetophenone	8.769	105	363298	38.050	ng	96
24) Nitrobenzene	9.151	77	283463	42.314	ng	96
25) Isophorone	9.674	82	500172	35.785	ng	99
26) 2-Nitrophenol	9.856	139	124211	51.689	ng	# 85
27) 2,4-Dimethylphenol	9.921	122	204942	36.502	ng	98
28) bis(2-Chloroethoxy)met...	10.156	93	292669	35.467	ng	98
29) 2,4-Dichlorophenol	10.397	162	224291	43.503	ng	98
30) 1,2,4-Trichlorobenzene	10.608	180	244316	41.848	ng	98
31) Naphthalene	10.796	128	693761	37.038	ng	99
32) Benzoic acid	10.091	122	169664	45.786	ng	97
33) 4-Chloroaniline	10.902	127	311825	35.842	ng	98
34) Hexachlorobutadiene	11.084	225	181932	46.759	ng	98
35) Caprolactam	11.689	113	74813m	36.637	ng	
36) 4-Chloro-3-methylphenol	12.030	107	258766	39.288	ng	97
37) 2-Methylnaphthalene	12.406	142	515673	37.446	ng	98
38) 1-Methylnaphthalene	12.623	142	486364	37.386	ng	97
40) 1,2,4,5-Tetrachloroben...	12.770	216	327743	42.765	ng	99
41) Hexachlorocyclopentadiene	12.753	237	185510	47.605	ng	99
43) 2,4,6-Trichlorophenol	13.011	196	217818	44.139	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.088	196	256349	43.529	ng	96
46) 1,1'-Biphenyl	13.417	154	695646	36.945	ng	98
47) 2-Chloronaphthalene	13.458	162	532178	37.200	ng	98
48) 2-Nitroaniline	13.652	65	213309	46.288	ng	97
49) Acenaphthylene	14.304	152	881868	36.687	ng	100
50) Dimethylphthalate	14.039	163	770371	36.562	ng	99
51) 2,6-Dinitrotoluene	14.151	165	162658	43.471	ng	94
52) Acenaphthene	14.645	154	582861m	38.620	ng	
53) 3-Nitroaniline	14.480	138	172126	41.972	ng	# 98
54) 2,4-Dinitrophenol	14.686	184	94495	64.430	ng	90
55) Dibenzofuran	14.979	168	875680	36.784	ng	96
56) 4-Nitrophenol	14.791	139	133480	43.334	ng	96
57) 2,4-Dinitrotoluene	14.944	165	235916	48.084	ng	90
58) Fluorene	15.632	166	711591	37.245	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.209	232	237433	46.263	ng	95
60) Diethylphthalate	15.408	149	760937	35.896	ng	99
61) 4-Chlorophenyl-phenyle...	15.626	204	396062	39.184	ng	96
62) 4-Nitroaniline	15.649	138	179067	46.483	ng	94
63) Azobenzene	15.919	77	703429	34.053	ng	94
65) 4,6-Dinitro-2-methylph...	15.708	198	144008	54.723	ng	98
66) n-Nitrosodiphenylamine	15.843	169	667901	37.426	ng	97
67) 4-Bromophenyl-phenylether	16.519	248	279265	44.063	ng	98
68) Hexachlorobenzene	16.636	284	301480	42.138	ng	96
69) Atrazine	16.795	200	232745	37.295	ng	97
70) Pentachlorophenol	16.977	266	215246	46.107	ng	97
71) Phenanthrene	17.371	178	1187163	36.552	ng	99
72) Anthracene	17.465	178	1222673	37.068	ng	100
73) Carbazole	17.729	167	1052059	36.176	ng	99
74) Di-n-butylphthalate	18.305	149	1268342	34.982	ng	99
75) Fluoranthene	19.380	202	1454938	36.917	ng	96
77) Benzidine	19.562	184	519071	34.234	ng	99
78) Pyrene	19.744	202	1477139	36.680	ng	100
80) Butylbenzylphthalate	20.643	149	554969	37.778	ng	97
81) Benzo(a)anthracene	21.572	228	1492392	36.768	ng	99
82) 3,3'-Dichlorobenzidine	21.489	252	538351	39.281	ng	97
83) Chrysene	21.636	228	1413136	36.122	ng	97
84) Bis(2-ethylhexyl)phtha...	21.501	149	779048	33.272	ng	# 97
85) Di-n-octyl phthalate	22.694	149	1367389	35.847	ng	100
87) Indeno(1,2,3-cd)pyrene	28.293	276	1754392	36.917	ng	# 96
88) Benzo(b)fluoranthene	23.740	252	1406811	36.286	ng	# 98
89) Benzo(k)fluoranthene	23.810	252	1451057	36.567	ng	# 98
90) Benzo(a)pyrene	24.586	252	1389763	36.887	ng	97
91) Dibenzo(a,h)anthracene	28.364	278	1453075	36.272	ng	# 95
92) Benzo(g,h,i)perylene	29.404	276	1422842	36.511	ng	# 96

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

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