

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083024\
 Data File : BM047401.D
 Acq On : 30 Aug 2024 11:28
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 09/02/2024

Supervised By :mohammad ahmed 09/02/2024

Quant Time: Aug 30 12:06:48 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082624.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Aug 26 18:11:40 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.345	152	235590	20.000	ng	0.00
21) Naphthalene-d8	10.104	136	917193	20.000	ng	0.00
39) Acenaphthene-d10	14.010	164	653125	20.000	ng	0.00
64) Phenanthrene-d10	16.780	188	1437118	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1280997	20.000	ng	0.00
86) Perylene-d12	23.727	264	1509652	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.004	112	988246	80.479	ng	0.00
7) Phenol-d6	6.575	99	1338566	80.020	ng	0.00
23) Nitrobenzene-d5	8.498	82	1326006	80.217	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	684410	86.856	ng	0.00
45) 2-Fluorobiphenyl	12.616	172	3398435	79.949	ng	0.00
79) Terphenyl-d14	19.444	244	5549418	77.110	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.016	88	190689	37.939	ng	99
3) Pyridine	3.393	79	545246	39.086	ng	99
4) n-Nitrosodimethylamine	3.316	42	213960	37.660	ng	98
6) Aniline	6.698	93	606398	40.025	ng	99
8) 2-Chlorophenol	6.928	128	550941	40.085	ng	97
9) Benzaldehyde	6.516	77	388056	47.138	ng	96
10) Phenol	6.598	94	654061	39.741	ng	99
11) bis(2-Chloroethyl)ether	6.798	93	547306	38.374	ng	99
12) 1,3-Dichlorobenzene	7.234	146	678453	40.368	ng	99
13) 1,4-Dichlorobenzene	7.381	146	682616	40.104	ng	99
14) 1,2-Dichlorobenzene	7.686	146	639630	39.653	ng	99
15) Benzyl Alcohol	7.604	79	523734	45.160	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.875	45	654823m	38.291	ng	
17) 2-Methylphenol	7.816	107	468625	40.983	ng	97
18) Hexachloroethane	8.392	117	257857	40.525	ng	99
19) n-Nitroso-di-n-propyla...	8.157	70	433813	37.726	ng	98
20) 3+4-Methylphenols	8.145	107	643942	40.504	ng	98
22) Acetophenone	8.169	105	851392	39.997	ng	# 99
24) Nitrobenzene	8.539	77	682215	40.128	ng	98
25) Isophorone	9.063	82	1261174	40.368	ng	99
26) 2-Nitrophenol	9.239	139	360394	43.096	ng	98
27) 2,4-Dimethylphenol	9.327	122	402167	42.606	ng	99
28) bis(2-Chloroethoxy)met...	9.545	93	769181	40.205	ng	98
29) 2,4-Dichlorophenol	9.786	162	637880	43.563	ng	98
30) 1,2,4-Trichlorobenzene	9.975	180	704204	41.637	ng	100
31) Naphthalene	10.151	128	1813905	40.491	ng	99
32) Benzoic acid	9.533	122	471006	41.788	ng	97
33) 4-Chloroaniline	10.286	127	707114	42.452	ng	99
34) Hexachlorobutadiene	10.427	225	448832	42.493	ng	99
35) Caprolactam	11.104	113	215806	45.496	ng	98
36) 4-Chloro-3-methylphenol	11.451	107	639103	42.980	ng	99
37) 2-Methylnaphthalene	11.780	142	1336174	40.907	ng	99
38) 1-Methylnaphthalene	12.004	142	1317738	40.791	ng	99
40) 1,2,4,5-Tetrachloroben...	12.163	216	792820	40.993	ng	99
41) Hexachlorocyclopentadiene	12.127	237	231223	37.881	ng	98
43) 2,4,6-Trichlorophenol	12.433	196	549348	41.435	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.521	196	622160	42.255	ng	98
46) 1,1'-Biphenyl	12.827	154	1818932	39.885	ng	99
47) 2-Chloronaphthalene	12.868	162	1418002	39.677	ng	99
48) 2-Nitroaniline	13.104	65	420716	42.026	ng	95
49) Acenaphthylene	13.727	152	2108034	40.585	ng	99
50) Dimethylphthalate	13.492	163	1893257	42.070	ng	99
51) 2,6-Dinitrotoluene	13.615	165	457486	43.662	ng	96
52) Acenaphthene	14.074	154	1336809	40.542	ng	99
53) 3-Nitroaniline	13.951	138	430482	44.523	ng	100
54) 2,4-Dinitrophenol	14.186	184	286304	44.370	ng	94
55) Dibenzofuran	14.415	168	2174004	40.933	ng	99
56) 4-Nitrophenol	14.327	139	317028	41.947	ng	98
57) 2,4-Dinitrotoluene	14.427	165	601285	44.330	ng	# 88
58) Fluorene	15.074	166	1774869	41.292	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.668	232	504016	41.675	ng	100
60) Diethylphthalate	14.874	149	1871557	42.160	ng	99
61) 4-Chlorophenyl-phenyle...	15.074	204	906557	41.460	ng	98
62) 4-Nitroaniline	15.133	138	438028	48.974	ng	97
63) Azobenzene	15.374	77	1549934	40.095	ng	97
65) 4,6-Dinitro-2-methylph...	15.204	198	387949	43.982	ng	97
66) n-Nitrosodiphenylamine	15.304	169	1540714	39.180	ng	99
67) 4-Bromophenyl-phenylether	15.974	248	605150	39.915	ng	96
68) Hexachlorobenzene	16.086	284	706233	40.331	ng	98
69) Atrazine	16.274	200	523804	40.752	ng	99
70) Pentachlorophenol	16.451	266	443479	39.172	ng	98
71) Phenanthrene	16.821	178	2817514	40.370	ng	100
72) Anthracene	16.915	178	2800515	41.162	ng	99
73) Carbazole	17.198	167	2830916	43.360	ng	100
74) Di-n-butylphthalate	17.774	149	3384167	42.568	ng	99
75) Fluoranthene	18.862	202	3583659	43.730	ng	99
77) Benzidine	19.068	184	1521856	71.172	ng	99
78) Pyrene	19.227	202	3706717	37.728	ng	100
80) Butylbenzylphthalate	20.156	149	1582441	41.159	ng	98
81) Benzo(a)anthracene	21.009	228	3417519	40.910	ng	99
82) 3,3'-Dichlorobenzidine	20.950	252	1266742	45.078	ng	99
83) Chrysene	21.068	228	3194421	40.651	ng	100
84) Bis(2-ethylhexyl)phtha...	20.956	149	2158255	39.718	ng	99
85) Di-n-octyl phthalate	21.985	149	3841636	42.003	ng	99
87) Indeno(1,2,3-cd)pyrene	26.756	276	3850688	40.746	ng	99
88) Benzo(b)fluoranthene	22.885	252	3564211	40.467	ng	100
89) Benzo(k)fluoranthene	22.944	252	3276886	39.247	ng	100
90) Benzo(a)pyrene	23.603	252	3100626	40.708	ng	100
91) Dibenzo(a,h)anthracene	26.785	278	3098503	40.472	ng	100
92) Benzo(g,h,i)perylene	27.685	276	3223872	40.505	ng	99

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

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