

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102123\
 Data File : BM042390.D
 Acq On : 20 Oct 2023 22:42
 Operator : MA/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDICC080

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 10/23/2023

Supervised By :mohammad ahmed 10/23/2023

Quant Time: Oct 21 02:28:28 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102123.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Oct 21 02:15:02 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	118108	20.000	ng	0.00
21) Naphthalene-d8	10.822	136	496219	20.000	ng	0.00
39) Acenaphthene-d10	14.645	164	268130	20.000	ng	0.00
64) Phenanthrene-d10	17.398	188	538430	20.000	ng	0.00
76) Chrysene-d12	21.580	240	500239	20.000	ng	0.00
86) Perylene-d12	24.050	264	501660	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	1379706	165.100	ng	0.00
7) Phenol-d6	7.151	99	1909436	170.498	ng	0.00
23) Nitrobenzene-d5	9.204	82	1928727	171.875	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0d	0.000	ng	
45) 2-Fluorobiphenyl	13.269	172	3442124	172.871	ng	0.00
79) Terphenyl-d14	20.009	244	5058724	179.346	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.393	88	310713	78.960	ng	99
3) Pyridine	3.810	79	959263m	83.067	ng	
4) n-Nitrosodimethylamine	3.740	42	478018	82.005	ng	95
6) Aniline	7.345	93	1236782	86.739	ng	99
8) 2-Chlorophenol	7.557	128	704158	84.917	ng	97
9) Benzaldehyde	7.157	77	434819	66.661	ng	99
10) Phenol	7.181	94	982996	84.946	ng	99
11) bis(2-Chloroethyl)ether	7.440	93	793413	83.510	ng	99
12) 1,3-Dichlorobenzene	7.881	146	707394	82.476	ng	99
13) 1,4-Dichlorobenzene	8.034	146	719646	83.354	ng	100
14) 1,2-Dichlorobenzene	8.351	146	686239	82.777	ng	99
15) Benzyl Alcohol	8.257	79	745429	89.791	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.528	45	1249781	83.137	ng	99
17) 2-Methylphenol	8.439	107	662982	86.821	ng	96
18) Hexachloroethane	9.063	117	279981	84.237	ng	97
19) n-Nitroso-di-n-propyla...	8.822	70	687656	90.594	ng	98
20) 3+4-Methylphenols	8.775	107	917034	89.533	ng	98
22) Acetophenone	8.851	105	1144803	84.565	ng	# 98
24) Nitrobenzene	9.245	77	964940	85.950	ng	99
25) Isophorone	9.757	82	1741158	91.301	ng	99
27) 2,4-Dimethylphenol	9.981	122	687789	88.914	ng	98
28) bis(2-Chloroethoxy)met...	10.239	93	1032601	85.463	ng	100
29) 2,4-Dichlorophenol	10.463	162	613223	91.158	ng	97
30) 1,2,4-Trichlorobenzene	10.675	180	627034	86.948	ng	98
31) Naphthalene	10.875	128	2167745	85.050	ng	99
33) 4-Chloroaniline	11.010	127	977729	88.354	ng	99
34) Hexachlorobutadiene	11.098	225	347978	89.170	ng	99
35) Caprolactam	11.851	113	234825	83.006	ng	95
36) 4-Chloro-3-methylphenol	12.104	107	742819	91.504	ng	100
37) 2-Methylnaphthalene	12.474	142	1500299	87.442	ng	99
38) 1-Methylnaphthalene	12.692	142	1404559	87.383	ng	99
40) 1,2,4,5-Tetrachloroben...	12.827	216	700694	90.394	ng	99
43) 2,4,6-Trichlorophenol	13.080	196	473645	94.391	ng	99
44) 2,4,5-Trichlorophenol	13.145	196	555340	93.621	ng	98
46) 1,1'-Biphenyl	13.480	154	1822647	85.687	ng	99
47) 2-Chloronaphthalene	13.533	162	1329688	85.206	ng	99

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48) 2-Nitroaniline	13.769	65	583392	81.839	ng	97
49) Acenaphthylene	14.374	152	2196229	87.173	ng	100
50) Dimethylphthalate	14.116	163	1706059	86.745	ng	99
51) 2,6-Dinitrotoluene	14.257	165	373107	81.852	ng	93
52) Acenaphthene	14.710	154	1318455	86.561	ng	99
53) 3-Nitroaniline	14.592	138	442890	81.653	ng	94
54) 2,4-Dinitrophenol	14.792	184	229908	83.214	ng	96
55) Dibenzofuran	15.045	168	2092145	86.843	ng	98
56) 4-Nitrophenol	14.874	139	363046	90.634	ng	99
57) 2,4-Dinitrotoluene	15.033	165	506164	82.071	ng	92
58) Fluorene	15.692	166	1679742	88.359	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.263	232	421792	94.654	ng	95
60) Diethylphthalate	15.457	149	1697920	87.484	ng	98
61) 4-Chlorophenyl-phenyle...	15.686	204	824322	92.074	ng	93
62) 4-Nitroaniline	15.762	138	425737	81.388	ng	98
63) Azobenzene	15.980	77	2043086	87.741	ng	97
65) 4,6-Dinitro-2-methylph...	15.786	198	300227	82.313	ng	92
66) n-Nitrosodiphenylamine	15.910	169	1422917	87.740	ng	99
67) 4-Bromophenyl-phenylether	16.580	248	475810	95.348	ng	96
68) Hexachlorobenzene	16.662	284	492227	92.279	ng	97
69) Atrazine	16.857	200	356446	86.672	ng	99
70) Pentachlorophenol	17.027	266	369699	82.796	ng	100
71) Phenanthrene	17.445	178	2554301	87.205	ng	100
72) Anthracene	17.533	178	2611753	90.032	ng	99
73) Carbazole	17.821	167	2413264	88.677	ng	99
74) Di-n-butylphthalate	18.339	149	2875855	81.372	ng	100
75) Fluoranthene	19.451	202	3027052	93.835	ng	98
77) Benzidine	19.656	184	589329	81.449	ng	100
78) Pyrene	19.821	202	3165929	90.568	ng	100
80) Butylbenzylphthalate	20.697	149	1329594	81.472	ng	96
81) Benzo(a)anthracene	21.562	228	3002430	89.531	ng	99
82) 3,3'-Dichlorobenzidine	21.503	252	934350	93.334	ng	98
83) Chrysene	21.621	228	2829221	87.839	ng	99
84) Bis(2-ethylhexyl)phtha...	21.444	149	1888107	81.304	ng	# 95
85) Di-n-octyl phthalate	22.391	149	3261593	80.844	ng	99
87) Indeno(1,2,3-cd)pyrene	26.650	276	3374289	92.040	ng	96
88) Benzo(b)fluoranthene	23.286	252	2784040	92.198	ng	98
89) Benzo(k)fluoranthene	23.338	252	2873279	91.651	ng	98
90) Benzo(a)pyrene	23.944	252	2724610	93.745	ng	97
91) Dibenzo(a,h)anthracene	26.668	278	2804188	91.998	ng	97
92) Benzo(g,h,i)perylene	27.462	276	2710161	91.554	ng	96

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

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