

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032423\
 Data File : BM039146.D
 Acq On : 24 Mar 2023 17:51
 Operator : CG/JU
 Sample : 02047-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 R-PITMSD

Quant Time: Mar 25 01:06:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 24 15:04:20 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/27/2023
 Supervised By : Jagrut Upadhyay 03/27/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	263680	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	991761	20.000	ng	0.00
39) Acenaphthene-d10	14.592	164	502155	20.000	ng	0.00
64) Phenanthrene-d10	17.339	188	900734	20.000	ng	0.00
76) Chrysene-d12	21.521	240	691877	20.000	ng	0.00
86) Perylene-d12	23.944	264	781653	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1226088	70.646	ng	0.00
7) Phenol-d6	7.116	99	1565100	69.427	ng	0.00
23) Nitrobenzene-d5	9.122	82	962227	47.658	ng	0.00
42) 2,4,6-Tribromophenol	16.086	330	387526	66.710	ng	0.00
45) 2-Fluorobiphenyl	13.222	172	1822397	47.244	ng	0.00
79) Terphenyl-d14	19.956	244	1822571	48.163	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.375	88	312599	40.486	ng	100
3) Pyridine	3.787	79	788272	36.812	ng	99
4) n-Nitrosodimethylamine	3.693	42	359618	41.270	ng	93
6) Aniline	7.275	93	972846	32.867	ng	99
8) 2-Chlorophenol	7.516	128	864656	46.796	ng	98
9) Benzaldehyde	7.087	77	583090	54.355	ng	97
10) Phenol	7.145	94	1094006	45.699	ng	97
11) bis(2-Chloroethyl)ether	7.375	93	858640	44.116	ng	98
12) 1,3-Dichlorobenzene	7.840	146	947903	48.274	ng	99
13) 1,4-Dichlorobenzene	7.987	146	963395	48.241	ng	99
14) 1,2-Dichlorobenzene	8.304	146	914264	47.527	ng	100
15) Benzyl Alcohol	8.198	79	735540m	45.522	ng	
16) 2,2'-oxybis(1-Chloropr...	8.487	45	1075481	43.752	ng	99
17) 2-Methylphenol	8.404	107	697983	42.444	ng	99
18) Hexachloroethane	9.034	117	363347	49.313	ng	96
19) n-Nitroso-di-n-propyla...	8.769	70	622429	43.836	ng	98
20) 3+4-Methylphenols	8.734	107	928065	42.266	ng	99
22) Acetophenone	8.781	105	1269798	45.635	ng	# 98
24) Nitrobenzene	9.163	77	924573	43.568	ng	97
25) Isophorone	9.692	82	1653506	44.504	ng	100
26) 2-Nitrophenol	9.875	139	450988	48.122	ng	96
27) 2,4-Dimethylphenol	9.939	122	754386	46.599	ng	99
28) bis(2-Chloroethoxy)met...	10.175	93	1040285	43.576	ng	100
29) 2,4-Dichlorophenol	10.416	162	722191	47.053	ng	99
30) 1,2,4-Trichlorobenzene	10.628	180	791633	47.199	ng	99
31) Naphthalene	10.816	128	2492274	44.070	ng	100
32) Benzoic acid	10.086	122	547408	46.315	ng	97
33) 4-Chloroaniline	10.928	127	258557	10.728	ng	99
34) Hexachlorobutadiene	11.098	225	456213	47.432	ng	97
35) Caprolactam	11.728	113	214139	45.841	ng	100
36) 4-Chloro-3-methylphenol	12.057	107	724653	44.338	ng	99
37) 2-Methylnaphthalene	12.422	142	1605887	42.367	ng	99
38) 1-Methylnaphthalene	12.639	142	1500433	42.241	ng	99
40) 1,2,4,5-Tetrachloroben...	12.792	216	797535	48.854	ng	99
41) Hexachlorocyclopentadiene	12.769	237	1087094	121.316	ng	99
43) 2,4,6-Trichlorophenol	13.033	196	522666	49.442	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.104	196	549792	44.433	ng	96
46) 1,1'-Biphenyl	13.433	154	2000820	47.032	ng	99
47) 2-Chloronaphthalene	13.474	162	1519076	47.354	ng	99
48) 2-Nitroaniline	13.680	65	450246	44.029	ng	94
49) Acenaphthylene	14.316	152	2307730	45.163	ng	99
50) Dimethylphthalate	14.057	163	1688202	43.690	ng	99
51) 2,6-Dinitrotoluene	14.174	165	373465	43.831	ng	97
52) Acenaphthene	14.657	154	1378645	43.902	ng	98
53) 3-Nitroaniline	14.504	138	239019	24.724	ng	96
54) 2,4-Dinitrophenol	14.710	184	474217	95.474	ng	91
55) Dibenzofuran	14.992	168	2127334	43.291	ng	99
56) 4-Nitrophenol	14.816	139	698572	91.030	ng	96
57) 2,4-Dinitrotoluene	14.963	165	490784	43.829	ng	95
58) Fluorene	15.645	166	1639169	42.583	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.221	232	436483	45.740	ng	97
60) Diethylphthalate	15.416	149	1612316	42.884	ng	99
61) 4-Chlorophenyl-phenyle...	15.633	204	799380	42.272	ng	98
62) 4-Nitroaniline	15.668	138	390988	38.800	ng	96
63) Azobenzene	15.927	77	1730344	43.495	ng	97
65) 4,6-Dinitro-2-methylph...	15.721	198	276250	46.810	ng	95
66) n-Nitrosodiphenylamine	15.851	169	1383011	45.545	ng	99
67) 4-Bromophenyl-phenylether	16.527	248	461332	47.419	ng	98
68) Hexachlorobenzene	16.645	284	522498	48.042	ng	98
69) Atrazine	16.804	200	428323	51.641	ng	99
70) Pentachlorophenol	16.992	266	662489	82.819	ng	99
71) Phenanthrene	17.380	178	2354623	44.541	ng	100
72) Anthracene	17.474	178	2360286	44.771	ng	100
73) Carbazole	17.745	167	2142178	44.522	ng	100
74) Di-n-butylphthalate	18.309	149	2570491	43.959	ng	99
75) Fluoranthene	19.398	202	2455998	44.001	ng	99
77) Benzidine	19.586	184	1216108	74.154	ng	98
78) Pyrene	19.756	202	2546113	48.996	ng	100
80) Butylbenzylphthalate	20.656	149	1039661	47.117	ng	97
81) Benzo(a)anthracene	21.503	228	2262257	45.404	ng	99
82) 3,3'-Dichlorobenzidine	21.439	252	615054	40.165	ng	99
83) Chrysene	21.556	228	2175911	44.902	ng	100
84) Bis(2-ethylhexyl)phtha...	21.427	149	1477010	46.324	ng	100
85) Di-n-octyl phthalate	22.362	149	2566086	48.369	ng	98
87) Indeno(1,2,3-cd)pyrene	26.480	276	3133044	52.676	ng	100
88) Benzo(b)fluoranthene	23.209	252	2330784	46.674	ng	100
89) Benzo(k)fluoranthene	23.256	252	2234193	42.990	ng	100
90) Benzo(a)pyrene	23.839	252	2084182	43.366	ng	100
91) Dibenzo(a,h)anthracene	26.497	278	2609218	51.800	ng	99
92) Benzo(g,h,i)perylene	27.256	276	2656391	53.903	ng	99

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

