

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032624\
 Data File : BM044772.D
 Acq On : 26 Mar 2024 17:13
 Operator : MA/JU
 Sample : P1864-01MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VNJ229-72-11976MS

Quant Time: Mar 26 17:41:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 02:28:24 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 03/27/2024
 Supervised By :mohammad ahmed 03/28/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.686	152	141515	20.000	ng	-0.02
21) Naphthalene-d8	10.469	136	555603	20.000	ng	-0.02
39) Acenaphthene-d10	14.333	164	322882	20.000	ng	-0.02
64) Phenanthrene-d10	17.092	188	655698	20.000	ng	-0.01
76) Chrysene-d12	21.291	240	594767	20.000	ng	-0.01
86) Perylene-d12	23.538	264	708604	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.298	112	867871	88.052	ng	0.00
7) Phenol-d6	6.869	99	1089940	85.956	ng	-0.01
23) Nitrobenzene-d5	8.851	82	663086	57.495	ng	-0.01
42) 2,4,6-Tribromophenol	15.839	330	375429	105.420	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	1324556	56.497	ng	-0.02
79) Terphenyl-d14	19.733	244	2022258	59.479	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.281	88	148039	35.734	ng	95
3) Pyridine	3.663	79	395851	34.552	ng	91
4) n-Nitrosodimethylamine	3.587	42	222084	48.370	ng	79
6) Aniline	7.028	93	270045	17.981	ng	98
8) 2-Chlorophenol	7.257	128	440570	44.877	ng	98
9) Benzaldehyde	6.845	77	70046m	10.547	ng	
10) Phenol	6.898	94	544837	43.023	ng	95
11) bis(2-Chloroethyl)ether	7.128	93	412040	39.850	ng	98
12) 1,3-Dichlorobenzene	7.575	146	467629	44.300	ng	98
13) 1,4-Dichlorobenzene	7.722	146	473697	44.525	ng	99
14) 1,2-Dichlorobenzene	8.033	146	452479	44.597	ng	99
15) Benzyl Alcohol	7.933	79	373446	44.796	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.204	45	617532m	43.077	ng	
17) 2-Methylphenol	8.133	107	349199	40.432	ng	98
18) Hexachloroethane	8.745	117	181532	46.155	ng	98
19) n-Nitroso-di-n-propyla...	8.498	70	306316	39.556	ng	94
20) 3+4-Methylphenols	8.463	107	482699	41.415	ng	98
22) Acetophenone	8.516	105	612130	41.936	ng	# 96
24) Nitrobenzene	8.892	77	471760	41.381	ng	99
25) Isophorone	9.416	82	841261	42.278	ng	97
26) 2-Nitrophenol	9.592	139	233857	47.538	ng	96
27) 2,4-Dimethylphenol	9.657	122	298522	34.507	ng	99
28) bis(2-Chloroethoxy)met...	9.898	93	515639	39.023	ng	99
29) 2,4-Dichlorophenol	10.122	162	382491	46.390	ng	98
30) 1,2,4-Trichlorobenzene	10.327	180	397142	45.276	ng	99
31) Naphthalene	10.516	128	1273825	41.776	ng	99
32) Benzoic acid	9.816	122	305790	44.570	ng	99
33) 4-Chloroaniline	10.639	127	219362	17.148	ng	100
34) Hexachlorobutadiene	10.780	225	228340	46.689	ng	98
35) Caprolactam	11.451	113	127856m	48.072	ng	
36) 4-Chloro-3-methylphenol	11.774	107	404027	44.656	ng	98
37) 2-Methylnaphthalene	12.133	142	844012	40.782	ng	97
38) 1-Methylnaphthalene	12.357	142	784290	40.662	ng	99
40) 1,2,4,5-Tetrachloroben...	12.504	216	403641	44.053	ng	99
41) Hexachlorocyclopentadiene	12.468	237	218374	41.835	ng	99
43) 2,4,6-Trichlorophenol	12.757	196	287032	46.393	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.833	196	312492	42.626	ng	98
46) 1,1'-Biphenyl	13.163	154	1058315	41.908	ng	99
47) 2-Chloronaphthalene	13.204	162	823448	44.096	ng	100
48) 2-Nitroaniline	13.427	65	290313	45.819	ng	99
49) Acenaphthylene	14.057	152	1380035	45.366	ng	100
50) Dimethylphthalate	13.810	163	983158	40.709	ng	99
51) 2,6-Dinitrotoluene	13.933	165	221372	44.017	ng	98
52) Acenaphthene	14.398	154	772578	42.013	ng	99
53) 3-Nitroaniline	14.262	138	147244	25.003	ng	97
54) 2,4-Dinitrophenol	14.474	184	165029	51.837	ng	95
55) Dibenzofuran	14.739	168	1234870	42.311	ng	98
56) 4-Nitrophenol	14.580	139	463809	95.896	ng	93
57) 2,4-Dinitrotoluene	14.727	165	315253	47.974	ng	# 90
58) Fluorene	15.392	166	995127	43.784	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.968	232	262910	46.191	ng	99
60) Diethylphthalate	15.174	149	1042413	42.638	ng	100
61) 4-Chlorophenyl-phenyle...	15.392	204	462803	42.416	ng	99
62) 4-Nitroaniline	15.433	138	249705	42.447	ng	93
63) Azobenzene	15.686	77	1086812	43.005	ng	95
65) 4,6-Dinitro-2-methylph...	15.492	198	118770	28.883	ng	97
66) n-Nitrosodiphenylamine	15.609	169	845603	41.379	ng	99
67) 4-Bromophenyl-phenylether	16.286	248	287230	42.350	ng	97
68) Hexachlorobenzene	16.392	284	342397	47.492	ng	94
69) Atrazine	16.574	200	294534	46.124	ng	95
70) Pentachlorophenol	16.745	266	482855	93.109	ng	99
71) Phenanthrene	17.133	178	1560092	43.054	ng	100
72) Anthracene	17.227	178	1608669	44.265	ng	99
73) Carbazole	17.503	167	1544538	45.125	ng	99
74) Di-n-butylphthalate	18.074	149	1923294	43.608	ng	99
75) Fluoranthene	19.156	202	1846833	45.594	ng	98
77) Benzidine	19.362	184	1324420	86.159	ng	99
78) Pyrene	19.521	202	1953536	44.814	ng	100
80) Butylbenzylphthalate	20.439	149	940857	47.891	ng	96
81) Benzo(a)anthracene	21.280	228	1887587	45.612	ng	100
82) 3,3'-Dichlorobenzidine	21.221	252	668498	46.903	ng	98
83) Chrysene	21.333	228	1768223	44.714	ng	99
84) Bis(2-ethylhexyl)phtha...	21.221	149	1310912	44.291	ng	99
85) Di-n-octyl phthalate	22.109	149	2362466	46.017	ng	99
87) Indeno(1,2,3-cd)pyrene	25.821	276	2333432	45.552	ng	97
88) Benzo(b)fluoranthene	22.868	252	1966342	47.219	ng	99
89) Benzo(k)fluoranthene	22.915	252	1807232	42.215	ng	98
90) Benzo(a)pyrene	23.444	252	1811669	44.604	ng	98
91) Dibenzo(a,h)anthracene	25.832	278	1906607	44.149	ng	97
92) Benzo(g,h,i)perylene	26.509	276	1827559	42.990	ng	95

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

