

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041521\
 Data File : BP005306.D
 Acq On : 15 Apr 2021 19:34
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 4/16/2021 11:32:36 AM

Quant Time: Apr 16 01:00:46 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 00:51:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	38764	20.00	ng	0.00
21) Naphthalene-d8	10.458	136	105187	20.00	ng	0.00
39) Acenaphthene-d10	14.328	164	71259	20.00	ng	0.00
64) Phenanthrene-d10	17.092	188	156982	20.00	ng	0.00
76) Chrysene-d12	21.198	240	191491	20.00	ng	0.00
86) Perylene-d12	23.469	264	179029	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.264	112	508768	137.74	ng	0.00
7) Phenol-d6	6.858	99	624164	148.10	ng	0.00
23) Nitrobenzene-d5	8.834	82	566185	169.74	ng	0.00
42) 2,4,6-Tribromophenol	15.840	330	177645	197.49	ng	0.00
45) 2-Fluorobiphenyl	12.946	172	952589	180.47	ng	0.00
79) Terphenyl-d14	19.675	244	1786060	183.23	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	109099	50.02	ng	98
3) Pyridine	3.593	79	280059m	57.28	ng	
4) n-Nitrosodimethylamine	3.511	42	81677	56.59	ng	# 65
6) Aniline	7.005	93	315090	69.46	ng	100
8) 2-Chlorophenol	7.234	128	222678	84.01	ng	98
9) Benzaldehyde	6.817	77	121270	43.70	ng	97
10) Phenol	6.887	94	311238	72.11	ng	99
11) bis(2-Chloroethyl)ether	7.105	93	227982	73.70	ng	99
12) 1,3-Dichlorobenzene	7.552	146	243196	84.31	ng	97
13) 1,4-Dichlorobenzene	7.699	146	241048	84.64	ng	100
14) 1,2-Dichlorobenzene	8.017	146	226229	83.57	ng	98
15) Benzyl Alcohol	7.917	79	177286	60.73	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.199	45	145716	80.90	ng	98
17) 2-Methylphenol	8.122	107	167411	70.71	ng	98
18) Hexachloroethane	8.734	117	92981	75.26	ng	95
19) n-Nitroso-di-n-propyla...	8.481	70	173136	66.79	ng	99
20) 3+4-Methylphenols	8.458	107	219247	70.70	ng	96
22) Acetophenone	8.493	105	343573	92.71	ng	# 97
24) Nitrobenzene	8.875	77	281286	81.58	ng	98
26) 2-Nitrophenol	9.581	139	82694	86.35	ng	94
27) 2,4-Dimethylphenol	9.646	122	134624	86.81	ng	99
28) bis(2-Chloroethoxy)met...	9.881	93	217856	82.59	ng	99
29) 2,4-Dichlorophenol	10.116	162	159077	92.30	ng	97
30) 1,2,4-Trichlorobenzene	10.322	180	188087	95.19	ng	99
31) Naphthalene	10.505	128	527600	91.61	ng	99
32) Benzoic acid	9.863	122	78486	77.15	ng	90
33) 4-Chloroaniline	10.628	127	194196	90.48	ng	98
34) Hexachlorobutadiene	10.781	225	136869	99.97	ng	97
35) Caprolactam	11.457	113	46044	80.49	ng	90
37) 2-Methylnaphthalene	12.128	142	363678	93.13	ng	98
38) 1-Methylnaphthalene	12.352	142	346994	93.89	ng	99
40) 1,2,4,5-Tetrachloroben...	12.504	216	218272	90.91	ng	99
43) 2,4,6-Trichlorophenol	12.757	196	132260	87.68	ng	99
44) 2,4,5-Trichlorophenol	12.834	196	143839	87.39	ng	99
46) 1,1'-Biphenyl	13.157	154	468493	86.74	ng	99
47) 2-Chloronaphthalene	13.199	162	376400	88.18	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
49) Acenaphthylene	14.051	152	573304	86.83	ng	100
50) Dimethylphthalate	13.799	163	451444	86.30	ng	98
51) 2,6-Dinitrotoluene	13.922	165	108155	92.97	ng	98
52) Acenaphthene	14.399	154	355398	88.12	ng	97
54) 2,4-Dinitrophenol	14.469	184	56763	89.19	ng	94
55) Dibenzofuran	14.734	168	617119	91.86	ng	97
56) 4-Nitrophenol	14.587	139	75743	85.00	ng	97
57) 2,4-Dinitrotoluene	14.722	165	158552	98.04	ng	96
58) Fluorene	15.387	166	507018	93.12	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.969	232	132571m	90.45	ng	
60) Diethylphthalate	15.169	149	459661	85.91	ng	99
61) 4-Chlorophenyl-phenyle...	15.387	204	271516	95.05	ng	98
62) 4-Nitroaniline	15.434	138	93994	86.35	ng	94
63) Azobenzene	15.681	77	637644	77.23	ng	99
65) 4,6-Dinitro-2-methylph...	15.493	198	90210	85.58	ng	97
66) n-Nitrosodiphenylamine	15.604	169	418875	88.31	ng	98
67) 4-Bromophenyl-phenylether	16.281	248	165970	92.54	ng	96
68) Hexachlorobenzene	16.398	284	186239	97.15	ng	97
69) Atrazine	16.569	200	138976	83.74	ng	95
70) Pentachlorophenol	16.751	266	101086	87.39	ng	93
71) Phenanthrene	17.140	178	787982	88.80	ng	99
72) Anthracene	17.228	178	774668	90.10	ng	99
73) Carbazole	17.510	167	724611	88.26	ng	98
75) Fluoranthene	19.122	202	982576	90.27	ng	99
77) Benzidine	19.310	184	207452	59.20	ng	99
78) Pyrene	19.475	202	1018537	85.58	ng	100
80) Butylbenzylphthalate	20.351	149	260878	58.66	ng	99
81) Benzo(a)anthracene	21.186	228	1045730	82.77	ng	100
82) 3,3'-Dichlorobenzidine	21.122	252	285538	77.25	ng	98
83) Chrysene	21.239	228	1033359	84.08	ng	100
84) Bis(2-ethylhexyl)phtha...	21.110	149	446100	63.84	ng	97
85) Di-n-octyl phthalate	21.986	149	796101	63.43	ng	99
87) Indeno(1,2,3-cd)pyrene	25.810	276	1141971	83.44	ng	99
88) Benzo(b)fluoranthene	22.786	252	1045661	89.23	ng	99
89) Benzo(k)fluoranthene	22.833	252	1041565	92.73	ng	99
90) Benzo(a)pyrene	23.375	252	951980	87.78	ng	98
91) Dibenzo(a,h)anthracene	25.821	278	996177	83.49	ng	99
92) Benzo(g,h,i)perylene	26.527	276	916778	79.46	ng	97

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

