

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
 Data File : BP005279.D
 Acq On : 13 Apr 2021 10:26
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad

4/14/2021 11:50:52 AM

Quant Time: Apr 14 02:38:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	32874	20.00	ng	# 0.00
21) Naphthalene-d8	10.463	136	93346	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	50434	20.00	ng	0.00
64) Phenanthrene-d10	17.098	188	100999	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	116730	20.00	ng	0.00
86) Perylene-d12	23.474	264	119286	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	270598	86.39	ng	0.00
7) Phenol-d6	6.863	99	283324	79.27	ng	0.00
23) Nitrobenzene-d5	8.840	82	250603	84.66	ng	0.00
42) 2,4,6-Tribromophenol	15.839	330	50443	79.23	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	327413	87.64	ng	0.00
79) Terphenyl-d14	19.680	244	488937	82.28	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.211	88	69683	37.68	ng	# 90
3) Pyridine	3.611	79	160714	38.76	ng	93
4) n-Nitrosodimethylamine	3.522	42	45218	36.94	ng	# 78
6) Aniline	7.010	93	148562	38.62	ng	# 85
8) 2-Chlorophenol	7.246	128	90839	40.41	ng	# 73
9) Benzaldehyde	6.828	77	88513	37.61	ng	# 72
10) Phenol	6.887	94	143003	39.07	ng	73
11) bis(2-Chloroethyl)ether	7.110	93	104626	39.88	ng	87
12) 1,3-Dichlorobenzene	7.563	146	98707	40.35	ng	# 83
13) 1,4-Dichlorobenzene	7.710	146	98594	40.82	ng	# 91
14) 1,2-Dichlorobenzene	8.028	146	89998	39.20	ng	# 89
15) Benzyl Alcohol	7.928	79	84158	33.99	ng	# 77
16) 2,2'-oxybis(1-Chloropr...	8.210	45	58573	38.35	ng	70
17) 2-Methylphenol	8.128	107	71614	35.66	ng	# 78
18) Hexachloroethane	8.740	117	41517	39.63	ng	# 85
19) n-Nitroso-di-n-propyla...	8.487	70	77354	35.18	ng	# 85
20) 3+4-Methylphenols	8.457	107	93626	35.60	ng	86
22) Acetophenone	8.504	105	137328	41.76	ng	# 88
24) Nitrobenzene	8.881	77	128667	42.05	ng	# 65
25) Isophorone	9.404	82	186663	37.80	ng	# 82
26) 2-Nitrophenol	9.587	139	32430	38.16	ng	# 29
27) 2,4-Dimethylphenol	9.657	122	54095	39.31	ng	# 77
28) bis(2-Chloroethoxy)met...	9.887	93	90636	38.72	ng	# 96
29) 2,4-Dichlorophenol	10.122	162	59607	38.97	ng	88
30) 1,2,4-Trichlorobenzene	10.328	180	72504	41.35	ng	96
31) Naphthalene	10.516	128	207174	40.53	ng	99
32) Benzoic acid	9.822	122	28957	32.07	ng	# 51
33) 4-Chloroaniline	10.634	127	72092	37.85	ng	# 73
34) Hexachlorobutadiene	10.793	225	53747	44.24	ng	99
35) Caprolactam	11.434	113	16338m	32.18	ng	
36) 4-Chloro-3-methylphenol	11.775	107	73062	36.93	ng	# 71
37) 2-Methylnaphthalene	12.140	142	135584	39.12	ng	# 90
38) 1-Methylnaphthalene	12.357	142	126999	38.72	ng	# 97
40) 1,2,4,5-Tetrachloroben...	12.510	216	76775	45.18	ng	# 96
41) Hexachlorocyclopentadiene	12.481	237	31526	40.26	ng	93
43) 2,4,6-Trichlorophenol	12.763	196	44714	41.88	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	12.840	196	47107	40.44	ng	#	83
46) 1,1'-Biphenyl	13.163	154	165593	43.32	ng		94
47) 2-Chloronaphthalene	13.204	162	131012	43.36	ng		93
48) 2-Nitroaniline	13.422	65	45451	35.56	ng	#	42
49) Acenaphthylene	14.057	152	195868	41.91	ng		99
50) Dimethylphthalate	13.804	163	145538	39.31	ng		99
51) 2,6-Dinitrotoluene	13.922	165	33202	40.32	ng	#	1
52) Acenaphthene	14.404	154	118423	41.48	ng		96
53) 3-Nitroaniline	14.257	138	28462	35.16	ng	#	7
54) 2,4-Dinitrophenol	14.475	184	16168	34.56	ng	#	50
55) Dibenzofuran	14.739	168	201910	42.47	ng		92
56) 4-Nitrophenol	14.592	139	22618	35.86	ng	#	1
57) 2,4-Dinitrotoluene	14.716	165	45354	39.62	ng	#	14
58) Fluorene	15.392	166	155848	40.44	ng		98
59) 2,3,4,6-Tetrachlorophenol	14.975	232	40418	38.96	ng	#	92
60) Diethylphthalate	15.175	149	145466	38.41	ng		99
61) 4-Chlorophenyl-phenyle...	15.392	204	83989	41.54	ng		96
62) 4-Nitroaniline	15.428	138	29365m	35.13	ng		
63) Azobenzene	15.686	77	243520	41.67	ng		77
65) 4,6-Dinitro-2-methylph...	15.492	198	24824	36.61	ng	#	70
66) n-Nitrosodiphenylamine	15.610	169	128441	42.09	ng		99
67) 4-Bromophenyl-phenylether	16.292	248	49929	43.27	ng		90
68) Hexachlorobenzene	16.404	284	54843	44.46	ng	#	91
69) Atrazine	16.575	200	41335	38.71	ng		98
70) Pentachlorophenol	16.757	266	28774	38.66	ng		93
71) Phenanthrene	17.145	178	238371	41.75	ng		100
72) Anthracene	17.233	178	231224	41.80	ng		98
73) Carbazole	17.516	167	214435	40.60	ng		99
74) Di-n-butylphthalate	18.069	149	227300	38.52	ng	#	94
75) Fluoranthene	19.127	202	278323	39.74	ng		98
77) Benzidine	19.316	184	74453	34.85	ng		99
78) Pyrene	19.480	202	290068	39.98	ng		98
80) Butylbenzylphthalate	20.357	149	87531	32.29	ng	#	60
81) Benzo(a)anthracene	21.186	228	296826	38.54	ng		99
82) 3,3'-Dichlorobenzidine	21.127	252	83841	37.21	ng		99
83) Chrysene	21.239	228	297806	39.75	ng		97
84) Bis(2-ethylhexyl)phtha...	21.116	149	145990	34.27	ng		99
85) Di-n-octyl phthalate	21.992	149	265398	34.69	ng	#	90
87) Indeno(1,2,3-cd)pyrene	25.809	276	379401	41.61	ng	#	93
88) Benzo(b)fluoranthene	22.786	252	315671	40.43	ng	#	98
89) Benzo(k)fluoranthene	22.833	252	299119	39.97	ng	#	98
90) Benzo(a)pyrene	23.374	252	290575	40.21	ng	#	97
91) Dibenzo(a,h)anthracene	25.815	278	328393	41.31	ng	#	95
92) Benzo(g,h,i)perylene	26.521	276	318382	41.41	ng	#	91

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

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