

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100720\
 Data File : BP003553.D
 Acq On : 07 Oct 2020 15:23
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 10/8/2020 12:29:49 PM

Quant Time: Oct 07 16:36:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 14:33:14 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.487	152	59731	20.00	ng	#-0.01
21) Naphthalene-d8	10.263	136	234112	20.00	ng	#-0.01
39) Acenaphthene-d10	14.145	164	143167	20.00	ng	-0.01
64) Phenanthrene-d10	16.910	188	309524	20.00	ng	# 0.00
76) Chrysene-d12	21.021	240	343222	20.00	ng	0.00
86) Perylene-d12	23.192	264	382760	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.105	112	277741	81.19	ng	0.00
7) Phenol-d6	6.699	99	392536	86.09	ng	0.00
23) Nitrobenzene-d5	8.640	82	365932	91.83	ng	-0.01
42) 2,4,6-Tribromophenol	15.657	330	157082	71.94	ng	0.00
45) 2-Fluorobiphenyl	12.763	172	792629	80.97	ng	-0.01
79) Terphenyl-d14	19.492	244	1308245	79.42	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.993	88	47243	33.61	ng	98
3) Pyridine	3.387	79	128986	34.46	ng	96
4) n-Nitrosodimethylamine	3.311	42	44380	39.64	ng	# 80
6) Aniline	6.822	93	226486	43.26	ng	90
8) 2-Chlorophenol	7.063	128	152868	40.12	ng	88
9) Benzaldehyde	6.640	77	100396	39.48	ng	83
10) Phenol	6.722	94	187711	43.10	ng	85
11) bis(2-Chloroethyl)ether	6.922	93	151532	43.91	ng	94
12) 1,3-Dichlorobenzene	7.381	146	179334	39.37	ng	# 93
13) 1,4-Dichlorobenzene	7.522	146	181065	39.30	ng	# 94
14) 1,2-Dichlorobenzene	7.840	146	173705	39.27	ng	95
15) Benzyl Alcohol	7.746	79	135504	44.72	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.028	45	120462	48.00	ng	89
17) 2-Methylphenol	7.963	107	130716	41.51	ng	# 92
18) Hexachloroethane	8.552	117	72835	42.42	ng	92
19) n-Nitroso-di-n-propyla...	8.299	70	119258	48.82	ng	# 93
20) 3+4-Methylphenols	8.293	107	176406	42.09	ng	# 87
22) Acetophenone	8.310	105	245090	42.63	ng	# 92
24) Nitrobenzene	8.687	77	185296	46.47	ng	# 84
25) Isophorone	9.204	82	322478	44.16	ng	# 90
26) 2-Nitrophenol	9.399	139	85471	39.25	ng	# 76
27) 2,4-Dimethylphenol	9.481	122	123163	39.99	ng	88
28) bis(2-Chloroethoxy)met...	9.693	93	211430	43.49	ng	97
29) 2,4-Dichlorophenol	9.951	162	142739	37.09	ng	93
30) 1,2,4-Trichlorobenzene	10.140	180	170574	37.36	ng	99
31) Naphthalene	10.310	128	492317	39.81	ng	99
32) Benzoic acid	9.693	122	31225m	20.62	ng	
33) 4-Chloroaniline	10.446	127	209668	39.91	ng	# 92
34) Hexachlorobutadiene	10.604	225	114137	36.62	ng	98
35) Caprolactam	11.210	113	48939	38.11	ng	# 76
36) 4-Chloro-3-methylphenol	11.604	107	164025	39.53	ng	87
37) 2-Methylnaphthalene	11.940	142	346616	38.69	ng	# 95
38) 1-Methylnaphthalene	12.163	142	328168m	38.58	ng	
40) 1,2,4,5-Tetrachloroben...	12.328	216	182735	38.98	ng	98
41) Hexachlorocyclopentadiene	12.304	237	50608	41.47	ng	100
43) 2,4,6-Trichlorophenol	12.587	196	110191	36.36	ng	99

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44) 2,4,5-Trichlorophenol	12.681	196	106238	34.08	ng	#	92
46) 1,1'-Biphenyl	12.975	154	438489	41.07	ng		98
47) 2-Chloronaphthalene	13.016	162	357261	40.40	ng		96
48) 2-Nitroaniline	13.234	65	109728	49.03	ng	#	68
49) Acenaphthylene	13.869	152	542066	40.12	ng		99
50) Dimethylphthalate	13.616	163	453745	39.69	ng		100
51) 2,6-Dinitrotoluene	13.740	165	97986	39.99	ng	#	70
52) Acenaphthene	14.210	154	330519	40.20	ng		100
53) 3-Nitroaniline	14.075	138	109000	41.05	ng	#	73
54) 2,4-Dinitrophenol	14.328	184	43914	41.69	ng	#	78
55) Dibenzofuran	14.551	168	514409	39.20	ng		93
56) 4-Nitrophenol	14.451	139	65961	40.64	ng	#	62
57) 2,4-Dinitrotoluene	14.551	165	135783	39.95	ng	#	48
58) Fluorene	15.210	166	416332	40.17	ng		99
59) 2,3,4,6-Tetrachlorophenol	14.804	232	104131m	35.52	ng		
60) Diethylphthalate	14.992	149	465522	39.92	ng		99
61) 4-Chlorophenyl-phenyle...	15.204	204	221416	39.13	ng		95
62) 4-Nitroaniline	15.251	138	109850	38.98	ng	#	64
63) Azobenzene	15.498	77	427053	48.04	ng		88
65) 4,6-Dinitro-2-methylph...	15.334	198	61007	37.35	ng		90
66) n-Nitrosodiphenylamine	15.422	169	369434	40.63	ng		99
67) 4-Bromophenyl-phenylether	16.098	248	144789	38.70	ng		93
68) Hexachlorobenzene	16.228	284	168973	38.02	ng		95
69) Atrazine	16.386	200	137414	38.53	ng		99
70) Pentachlorophenol	16.586	266	73329	41.95	ng		98
71) Phenanthrene	16.951	178	672451	39.93	ng		99
72) Anthracene	17.045	178	668876	40.30	ng		99
73) Carbazole	17.322	167	631924	40.09	ng		99
74) Di-n-butylphthalate	17.886	149	807568	40.07	ng	#	98
75) Fluoranthene	18.939	202	829184	39.16	ng		99
77) Benzidine	19.127	184	352923	33.16	ng		100
78) Pyrene	19.292	202	862387	40.29	ng		100
80) Butylbenzylphthalate	20.174	149	398604	40.70	ng	#	85
81) Benzo(a)anthracene	21.004	228	886378	39.95	ng		100
82) 3,3'-Dichlorobenzidine	20.939	252	331530	38.66	ng		99
83) Chrysene	21.057	228	863744	40.52	ng		100
84) Bis(2-ethylhexyl)phtha...	20.939	149	563500	40.87	ng		98
85) Di-n-octyl phthalate	21.786	149	1026305	40.74	ng	#	97
87) Indeno(1,2,3-cd)pyrene	25.386	276	1165690	40.18	ng		98
88) Benzo(b)fluoranthene	22.539	252	978497	40.48	ng		100
89) Benzo(k)fluoranthene	22.580	252	917089	39.59	ng		99
90) Benzo(a)pyrene	23.098	252	917515	40.26	ng		99
91) Dibenzo(a,h)anthracene	25.392	278	964364	40.28	ng		98
92) Benzo(g,h,i)perylene	26.057	276	940102	39.33	ng		98

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

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