

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040121\
 Data File : BP005130.D
 Acq On : 01 Apr 2021 11:40
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 4/2/2021 12:10:05 PM

Quant Time: Apr 01 12:24:29 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:23:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	33388	20.00	ng	# 0.00
21) Naphthalene-d8	10.499	136	132496	20.00	ng	# 0.00
39) Acenaphthene-d10	14.363	164	85183	20.00	ng	0.00
64) Phenanthrene-d10	17.128	188	168916	20.00	ng	# 0.00
76) Chrysene-d12	21.228	240	167664	20.00	ng	0.00
86) Perylene-d12	23.510	264	169255	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	179836	82.89	ng	0.00
7) Phenol-d6	6.887	99	259617	83.54	ng	0.00
23) Nitrobenzene-d5	8.869	82	239049	77.41	ng	0.00
42) 2,4,6-Tribromophenol	15.863	330	86192	77.66	ng	0.00
45) 2-Fluorobiphenyl	12.981	172	478172	75.94	ng	0.00
79) Terphenyl-d14	19.698	244	729233	79.18	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.246	88	37632	39.83	ng	97
3) Pyridine	3.646	79	99302	42.27	ng	95
4) n-Nitrosodimethylamine	3.558	42	34065	40.56	ng	# 90
6) Aniline	7.046	93	149192	42.36	ng	# 89
8) 2-Chlorophenol	7.275	128	96324	42.09	ng	85
9) Benzaldehyde	6.858	77	61251	38.70	ng	# 79
10) Phenol	6.917	94	131256	41.21	ng	87
11) bis(2-Chloroethyl)ether	7.140	93	95952	41.15	ng	92
12) 1,3-Dichlorobenzene	7.599	146	100929	39.70	ng	# 92
13) 1,4-Dichlorobenzene	7.746	146	102800	39.83	ng	# 94
14) 1,2-Dichlorobenzene	8.058	146	100507	40.59	ng	93
15) Benzyl Alcohol	7.952	79	100748	42.02	ng	# 87
16) 2,2'-oxybis(1-Chloropr...	8.234	45	78116	45.67	ng	94
17) 2-Methylphenol	8.152	107	86597	42.60	ng	# 92
18) Hexachloroethane	8.775	117	38979	39.63	ng	90
19) n-Nitroso-di-n-propyla...	8.522	70	84535	41.93	ng	# 93
20) 3+4-Methylphenols	8.487	107	118547	42.71	ng	89
22) Acetophenone	8.534	105	147726	38.78	ng	# 93
24) Nitrobenzene	8.911	77	124307	38.20	ng	# 79
25) Isophorone	9.440	82	231382	40.60	ng	# 89
26) 2-Nitrophenol	9.616	139	53715	42.56	ng	# 73
27) 2,4-Dimethylphenol	9.681	122	79443	39.14	ng	87
28) bis(2-Chloroethoxy)met...	9.916	93	117035	40.57	ng	98
29) 2,4-Dichlorophenol	10.152	162	89789	39.54	ng	93
30) 1,2,4-Trichlorobenzene	10.363	180	96924	38.00	ng	99
31) Naphthalene	10.552	128	293747	39.25	ng	99
32) Benzoic acid	9.846	122	68486	45.45	ng	# 77
33) 4-Chloroaniline	10.663	127	120594	39.82	ng	# 85
34) Hexachlorobutadiene	10.828	225	61846	37.68	ng	95
35) Caprolactam	11.469	113	32554	43.68	ng	86
36) 4-Chloro-3-methylphenol	11.799	107	108440	39.53	ng	85
37) 2-Methylnaphthalene	12.169	142	208713	38.62	ng	95
38) 1-Methylnaphthalene	12.387	142	193756m	38.25	ng	
40) 1,2,4,5-Tetrachloroben...	12.540	216	103766	37.41	ng	# 98
41) Hexachlorocyclopentadiene	12.516	237	59211	39.24	ng	100
43) 2,4,6-Trichlorophenol	12.787	196	76855	40.08	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	12.857	196	81460	39.23	ng	#	94
46) 1,1'-Biphenyl	13.193	154	250334	38.53	ng		97
47) 2-Chloronaphthalene	13.234	162	202712	38.33	ng		99
48) 2-Nitroaniline	13.446	65	69138	41.03	ng	#	61
49) Acenaphthylene	14.087	152	318500	38.72	ng		100
50) Dimethylphthalate	13.828	163	249028	37.96	ng		98
51) 2,6-Dinitrotoluene	13.946	165	61569	39.82	ng	#	60
52) Acenaphthene	14.428	154	194311	38.39	ng		99
53) 3-Nitroaniline	14.281	138	59182	41.58	ng	#	63
54) 2,4-Dinitrophenol	14.487	184	40392	43.14	ng	#	80
55) Dibenzofuran	14.769	168	312208	37.64	ng		98
56) 4-Nitrophenol	14.593	139	50931	43.63	ng	#	61
57) 2,4-Dinitrotoluene	14.740	165	82324	40.31	ng	#	63
58) Fluorene	15.422	166	255222	38.10	ng		99
59) 2,3,4,6-Tetrachlorophenol	14.998	232	73804	38.53	ng		95
60) Diethylphthalate	15.198	149	249024	38.80	ng		100
61) 4-Chlorophenyl-phenyle...	15.416	204	129884	36.54	ng		97
62) 4-Nitroaniline	15.451	138	62846	42.98	ng	#	74
63) Azobenzene	15.710	77	273466	37.28	ng		86
65) 4,6-Dinitro-2-methylph...	15.504	198	54865	44.51	ng		88
66) n-Nitrosodiphenylamine	15.634	169	221362	40.77	ng		97
67) 4-Bromophenyl-phenylether	16.316	248	79656	40.70	ng		97
68) Hexachlorobenzene	16.428	284	86174	39.15	ng		98
69) Atrazine	16.598	200	74991	41.58	ng		95
70) Pentachlorophenol	16.775	266	60918	41.92	ng		99
71) Phenanthrene	17.169	178	387962	39.62	ng		99
72) Anthracene	17.257	178	388190	40.57	ng		99
73) Carbazole	17.534	167	366448	40.77	ng		99
74) Di-n-butylphthalate	18.092	149	423396	42.80	ng	#	98
75) Fluoranthene	19.145	202	454965	39.49	ng		97
77) Benzidine	19.334	184	168228	45.73	ng		99
78) Pyrene	19.498	202	464340	40.30	ng		99
80) Butylbenzylphthalate	20.381	149	193082	45.16	ng	#	78
81) Benzo(a)anthracene	21.210	228	454575	39.91	ng		98
82) 3,3'-Dichlorobenzidine	21.145	252	164531	42.98	ng		100
83) Chrysene	21.263	228	442478	39.58	ng		97
84) Bis(2-ethylhexyl)phtha...	21.139	149	276841	44.17	ng		96
85) Di-n-octyl phthalate	22.022	149	476180	45.06	ng	#	94
87) Indeno(1,2,3-cd)pyrene	25.863	276	501324	39.50	ng		99
88) Benzo(b)fluoranthene	22.816	252	454271	37.80	ng		99
89) Benzo(k)fluoranthene	22.863	252	442953	38.46	ng		99
90) Benzo(a)pyrene	23.410	252	419002	38.88	ng		99
91) Dibenzo(a,h)anthracene	25.874	278	432938	39.45	ng		98
92) Benzo(g,h,i)perylene	26.580	276	422423	39.89	ng		97

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

