

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042921\
 Data File : BP005408.D
 Acq On : 29 Apr 2021 13:54
 Operator : CG/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

mohammad
 5/3/2021 10:48:04 AM

Quant Time: Apr 29 14:42:33 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP042921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 29 14:39:11 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.799	196	75926	44.17	ng	92
46) 1,1'-Biphenyl	13.116	154	201536	42.41	ng	97
47) 2-Chloronaphthalene	13.157	162	165076	42.19	ng	97
48) 2-Nitroaniline	13.381	65	55855	52.00	ng	94
49) Acenaphthylene	14.010	152	240687	42.51	ng	99
50) Dimethylphthalate	13.763	163	212506	43.85	ng	100
51) 2,6-Dinitrotoluene	13.887	165	47380	44.83	ng	93
52) Acenaphthene	14.357	154	148964	41.70	ng	97
53) 3-Nitroaniline	14.222	138	39130	44.78	ng	99
54) 2,4-Dinitrophenol	14.440	184	32017	53.16	ng	91
55) Dibenzofuran	14.698	168	268428	41.95	ng	95
56) 4-Nitrophenol	14.551	139	29507	42.42	ng	83
57) 2,4-Dinitrotoluene	14.687	165	69606	47.08	ng	94
58) Fluorene	15.351	166	226152	44.05	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.934	232	74035	41.39	ng	92
60) Diethylphthalate	15.134	149	211730	46.29	ng	97
61) 4-Chlorophenyl-phenyle...	15.351	204	140449	42.64	ng	98
62) 4-Nitroaniline	15.393	138	37396	45.61	ng	92
63) Azobenzene	15.645	77	261618	49.08	ng	96
65) 4,6-Dinitro-2-methylph...	15.457	198	50000	46.33	ng	92
66) n-Nitrosodiphenylamine	15.569	169	188431	42.01	ng	# 94
67) 4-Bromophenyl-phenylether	16.245	248	93489	41.68	ng	92
68) Hexachlorobenzene	16.357	284	106976	40.67	ng	# 90
69) Atrazine	16.534	200	85756	45.80	ng	99
70) Pentachlorophenol	16.716	266	59244	41.89	ng	97
71) Phenanthrene	17.098	178	358176	40.94	ng	99
72) Anthracene	17.192	178	359365	42.18	ng	99
73) Carbazole	17.475	167	319817	41.71	ng	99
74) Di-n-butylphthalate	18.034	149	358404	49.50	ng	# 96
75) Fluoranthene	19.086	202	493635	41.10	ng	98
77) Benzidine	19.281	184	167473	47.11	ng	100
78) Pyrene	19.439	202	509013	40.35	ng	99
80) Butylbenzylphthalate	20.322	149	172608	60.24	ng	90
81) Benzo(a)anthracene	21.151	228	575340	39.75	ng	99
82) 3,3'-Dichlorobenzidine	21.086	252	201250	44.73	ng	92
83) Chrysene	21.204	228	567846	39.38	ng	99
84) Bis(2-ethylhexyl)phtha...	21.075	149	273763	62.19	ng	98
85) Di-n-octyl phthalate	21.951	149	461381	53.97	ng	99
87) Indeno(1,2,3-cd)pyrene	25.727	276	798177	43.26	ng	98
88) Benzo(b)fluoranthene	22.739	252	628590	39.53	ng	99
89) Benzo(k)fluoranthene	22.780	252	623453m	39.94	ng	
90) Benzo(a)pyrene	23.322	252	583438	40.24	ng	99
91) Dibenzo(a,h)anthracene	25.733	278	693779	43.42	ng	100
92) Benzo(g,h,i)perylene	26.427	276	645522	43.13	ng	# 97

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

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