

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032422\
 Data File : BG052822.D
 Acq On : 24 Mar 2022 23:33
 Operator : CG/JU
 Sample : N2091-01MSD
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

MHBMSD

Manual Integrations

APPROVED

Reviewed By :Christian
 Giraldo

03/25/2022

Supervised By :Jagrut
 Upadhyay

03/25/2022

Quant Time: Mar 25 03:20:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 16:48:57 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.276	196	121929	57.523	ng	98
46) 1,1'-Biphenyl	13.599	154	327178	44.821	ng	96
47) 2-Chloronaphthalene	13.646	162	261068	45.414	ng	96
48) 2-Nitroaniline	13.834	65	105088	66.843	ng	90
49) Acenaphthylene	14.486	152	454862	49.637	ng	97
50) Dimethylphthalate	14.210	163	364596	46.999	ng	100
51) 2,6-Dinitrotoluene	14.328	165	82111	62.691	ng	# 80
52) Acenaphthene	14.827	154	334544	56.986	ng	99
53) 3-Nitroaniline	14.657	138	69612	45.420	ng	92
54) 2,4-Dinitrophenol	14.862	184	112731	157.900	ng	# 86
55) Dibenzofuran	15.162	168	445967	46.518	ng	98
56) 4-Nitrophenol	14.962	139	151271	120.998	ng	# 82
57) 2,4-Dinitrotoluene	15.115	165	119093	65.533	ng	# 91
58) Fluorene	15.808	166	369304	47.257	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.385	232	121339	53.556	ng	98
60) Diethylphthalate	15.573	149	391378	47.795	ng	100
61) 4-Chlorophenyl-phenyle...	15.796	204	202003	46.870	ng	98
62) 4-Nitroaniline	15.820	138	91078	56.240	ng	# 80
63) Azobenzene	16.090	77	404494	46.216	ng	97
65) 4,6-Dinitro-2-methylph...	15.878	198	73702	61.754	ng	93
66) n-Nitrosodiphenylamine	16.008	169	339516	45.048	ng	99
67) 4-Bromophenyl-phenylether	16.695	248	147710	45.766	ng	97
68) Hexachlorobenzene	16.818	284	164347	47.374	ng	95
69) Atrazine	16.953	200	148338	53.377	ng	98
70) Pentachlorophenol	17.159	266	217409	101.016	ng	95
71) Phenanthrene	17.553	178	657213	45.890	ng	98
72) Anthracene	17.641	178	666925	47.107	ng	99
73) Carbazole	17.905	167	611340	43.600	ng	100
74) Di-n-butylphthalate	18.463	149	703478	45.925	ng	99
75) Fluoranthene	19.556	202	834592	45.909	ng	99
77) Benzidine	19.726	184	446106	67.483	ng	98
78) Pyrene	19.914	202	825860	46.451	ng	100
80) Butylbenzylphthalate	20.795	149	311115	49.331	ng	97
81) Benzo(a)anthracene	21.776	228	807807	47.068	ng	99
82) 3,3'-Dichlorobenzidine	21.682	252	200748	33.123	ng	98
83) Chrysene	21.847	228	736110	45.623	ng	100
84) Bis(2-ethylhexyl)phtha...	21.676	149	424987	49.551	ng	99
85) Di-n-octyl phthalate	22.922	149	723943	50.702	ng	96
87) Indeno(1,2,3-cd)pyrene	28.908	276	891441	46.968	ng	# 96
88) Benzo(b)fluoranthene	24.061	252	845689	49.119	ng	96
89) Benzo(k)fluoranthene	24.132	252	787573	46.493	ng	97
90) Benzo(a)pyrene	24.960	252	805413	55.264	ng	100
91) Dibenzo(a,h)anthracene	28.978	278	771869	49.355	ng	98
92) Benzo(g,h,i)perylene	30.106	276	768898	49.689	ng	99

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

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