

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082321\
 Data File : BM031874.D
 Acq On : 23 Aug 2021 16:40
 Operator : CG/JU
 Sample : PB138616BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB138616BL

Quant Time: Aug 23 17:22:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 15:51:36 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.528	152	33717	20.000	ng	0.00
21) Naphthalene-d8	10.286	136	137300	20.000	ng	0.00
39) Acenaphthene-d10	14.168	164	84006	20.000	ng	0.00
64) Phenanthrene-d10	16.921	188	163814	20.000	ng	0.00
76) Chrysene-d12	21.127	240	117927	20.000	ng	0.00
86) Perylene-d12	23.274	264	106448	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.169	112	357604	156.024	ng	0.00
7) Phenol-d6	6.728	99	414283	134.574	ng	0.00
23) Nitrobenzene-d5	8.681	82	293885	87.161	ng	0.00
42) 2,4,6-Tribromophenol	15.668	330	117211	126.623	ng	0.00
45) 2-Fluorobiphenyl	12.780	172	570987	87.453	ng	0.00
79) Terphenyl-d14	19.574	244	740313	94.445	ng	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.175	88	68	0.091	ng	# 1
3) Pyridine	3.528	79	89	0.040	ng	# 1
4) n-Nitrosodimethylamine	3.416	42	161	0.164	ng	# 1
15) Benzyl Alcohol	7.522	79	1413	0.558	ng	# 50
24) Nitrobenzene	8.681	77	1290	0.374	ng	# 47
46) 1,1'-Biphenyl	12.780	154	1097	0.162	ng	# 1
52) Acenaphthene	14.168	154	87	0.017	ng	# 4
63) Azobenzene	15.674	77	1593	0.186	ng	# 32
66) n-Nitrosodiphenylamine	15.668	169	6199	1.121	ng	# 49
78) Pyrene	19.574	202	3569	0.378	ng	# 74
80) Butylbenzylphthalate	20.262	149	170	0.036	ng	# 19
81) Benzo(a)anthracene	21.127	228	272	0.031	ng	# 5
83) Chrysene	21.127	228	272	0.033	ng	# 7
84) Bis(2-ethylhexyl)phtha...	21.044	149	168	0.023	ng	# 55
85) Di-n-octyl phthalate	21.909	149	187	0.016	ng	# 1
87) Indeno(1,2,3-cd)pyrene	25.444	276	54	0.007	ng	# 51
88) Benzo(b)fluoranthene	22.650	252	228	0.029	ng	# 1
89) Benzo(k)fluoranthene	22.691	252	296	0.039	ng	# 1
90) Benzo(a)pyrene	23.197	252	55	0.008	ng	# 1

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

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