

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
 Data File : BF118968.D
 Acq On : 20 Feb 2020 19:41
 Operator : CG/JU
 Sample : L1599-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-BPM-03-011-ABCD

Quant Time: Feb 21 02:11:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 20 17:40:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	2485	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.05
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.79
64) Phenanthrene-d10	11.33	188	135	20.00	ng	0.05
76) Chrysene-d12	0.00	240	0	0.00	ng	-13.91
87) Perylene-d12	0.00	264	0	0.00	ng	-15.33

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	944599	6352.49	ng	0.00
7) Phenol-d6	6.40	99	1155684	6390.52	ng	0.00
23) Nitrobenzene-d5	7.32	82	754250	0.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.58	330	441759	0.00	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	1427806	0.00	ng	0.00
79) Terphenyl-d14	12.86	244	1754686	0.00	ng	0.00

Target Compounds

						Qvalue
9) Benzaldehyde	6.40	77	2075	17.174	ng	# 1
10) Phenol	6.41	94	11609	59.374	ng	# 96
15) Benzyl Alcohol	6.92	79	13492	103.227	ng	# 48
16) 2,2'-oxybis(1-Chloropropan	7.04	45	369	2.847	ng	# 52
19) n-Nitroso-di-n-propylamine	7.32	70	103448	981.689	ng	# 77
65) 4,6-Dinitro-2-methylphenol	10.57	198	930	1138.446	ng	# 1
66) n-Nitrosodiphenylamine	10.57	169	11593	2622.616	ng	# 41
73) Carbazole	11.51	167	1803	275.125	ng	# 58
74) Di-n-butylphthalate	11.94	149	231	29.107	ng	# 78
75) Fluoranthene	12.47	202	273	34.934	ng	# 64

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

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