

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070621\
 Data File : BN015348.D
 Acq On : 06 Jul 2021 21:02
 Operator : CG/JU
 Sample : M2898-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-03-6.5-11.5-062921

Quant Time: Jul 07 07:44:58 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 09:06:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.735	152	16286	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	76479	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	44261	0.40	ng	0.00
19) Phenanthrene-d10	17.090	188	90132	0.40	ng	0.00
29) Chrysene-d12	21.291	240	81553	0.40	ng	#-0.01
35) Perylene-d12	23.539	264	47570	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	5063	0.12	ng	0.00
5) Phenol-d6	6.905	99	3472	0.07	ng	0.00
8) Nitrobenzene-d5	8.873	82	17674	0.41	ng	0.00
11) 2-Methylnaphthalene-d10	12.087	152	40147	0.32	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	7189	0.60	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	55989	0.31	ng	0.00
27) Fluoranthene-d10	19.132	212	110527	0.42	ng	0.00
31) Terphenyl-d14	19.733	244	86433	0.45	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.738	266	1589	0.39	ng	# 35
32) Benzo(a)anthracene	21.280	228	14396	0.05	ng	97
33) Chrysene	21.323	228	13819	0.05	ng	98
34) Bis(2-ethylhexyl)phtha...	21.217	149	24729	0.40	ng	# 87
36) Indeno(1,2,3-cd)pyrene	25.798	276	5013	0.03	ng	# 90
37) Benzo(b)fluoranthene	22.865	252	10972	0.06	ng	# 77
38) Benzo(k)fluoranthene	22.909	252	10460	0.05	ng	# 84
39) Benzo(a)pyrene	23.443	252	8301	0.05	ng	# 65
40) Dibenzo(a,h)anthracene	25.809	278	4598	0.03	ng	# 81
41) Benzo(g,h,i)perylene	26.483	276	3624	0.02	ng	# 76

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

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