

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070621\
 Data File : BN015349.D
 Acq On : 06 Jul 2021 21:38
 Operator : CG/JU
 Sample : M2898-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EB-01-062921

Quant Time: Jul 07 01:33:15 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	15951	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	76022	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	44537	0.40	ng	0.00
19) Phenanthrene-d10	17.091	188	91106	0.40	ng	0.00
29) Chrysene-d12	21.291	240	81765	0.40	ng	#-0.01
35) Perylene-d12	23.539	264	46257	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	5504	0.14	ng	0.00
5) Phenol-d6	6.905	99	3850	0.08	ng	0.00
8) Nitrobenzene-d5	8.873	82	17789	0.42	ng	0.00
11) 2-Methylnaphthalene-d10	12.087	152	42337	0.34	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	6825	0.58	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	58913	0.33	ng	0.00
27) Fluoranthene-d10	19.133	212	112951	0.42	ng	0.00
31) Terphenyl-d14	19.733	244	88185	0.46	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.750	266	79	0.25	ng	# 79
32) Benzo(a)anthracene	21.281	228	14511	0.05	ng	97
33) Chrysene	21.323	228	13937	0.05	ng	98
34) Bis(2-ethylhexyl)phtha...	21.217	149	41424	0.56	ng	# 88
36) Indeno(1,2,3-cd)pyrene	25.792	276	4399	0.03	ng	# 89
37) Benzo(b)fluoranthene	22.865	252	10536	0.06	ng	# 71
38) Benzo(k)fluoranthene	22.910	252	10131	0.05	ng	# 81
39) Benzo(a)pyrene	23.440	252	7806	0.05	ng	# 58
40) Dibenzo(a,h)anthracene	25.809	278	4047	0.03	ng	# 78
41) Benzo(g,h,i)perylene	26.483	276	3067	0.02	ng	# 72

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

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