

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030322\
 Data File : BN018807.D
 Acq On : 03 Mar 2022 01:41
 Operator : CG/JU
 Sample : N1661-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-106

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/03/2022
 Supervised By : Jagrut Upadhyay 03/03/2022

Quant Time: Mar 03 04:35:00 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 03 00:44:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.912	152	5880	0.400	ng	0.00
7) Naphthalene-d8	10.712	136	15778	0.400	ng	#-0.01
13) Acenaphthene-d10	14.538	164	9767	0.400	ng	0.00
19) Phenanthrene-d10	17.277	188	22748	0.400	ng	0.00
29) Chrysene-d12	21.458	240	18120	0.400	ng	# 0.00
35) Perylene-d12	23.855	264	13666	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	1649	0.128	ng	0.00
5) Phenol-d6	7.067	99	1201	0.078	ng	0.00
8) Nitrobenzene-d5	9.067	82	3124m	0.272	ng	0.00
11) 2-Methylnaphthalene-d10	12.303	152	6916	0.269	ng	0.00
14) 2,4,6-Tribromophenol	16.025	330	1067	0.237	ng	0.00
15) 2-Fluorobiphenyl	13.169	172	12825	0.306	ng	0.00
27) Fluoranthene-d10	19.307	212	26440	0.395	ng	0.00
31) Terphenyl-d14	19.902	244	18154	0.469	ng	0.00
Target Compounds						
9) Naphthalene	10.765	128	3118	0.069	ng	# 93
12) 2-Methylnaphthalene	12.378	142	2564	0.081	ng	98
34) Bis(2-ethylhexyl)phtha...	21.369	149	2952	0.070	ng	99

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

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