

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041122\
 Data File : BN019409.D
 Acq On : 11 Apr 2022 13:12
 Operator : CG/JU
 Sample : N2334-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-18R

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/12/2022
 Supervised By : Jagrut Upadhyay 04/13/2022

Quant Time: Apr 12 07:31:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN040422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 04 17:40:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.826	152	5027	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	13174	0.400	ng	# 0.00
13) Acenaphthene-d10	14.463	164	7371	0.400	ng	0.00
19) Phenanthrene-d10	17.205	188	13862	0.400	ng	0.00
29) Chrysene-d12	21.387	240	10252	0.400	ng	# 0.00
35) Perylene-d12	23.738	264	8980	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.377	112	14	0.001	ng	0.00
5) Phenol-d6	7.002	99	45	0.003	ng	0.01
8) Nitrobenzene-d5	8.993	82	2643m	0.221	ng	0.01
11) 2-Methylnaphthalene-d10	12.223	152	5979	0.258	ng	0.00
14) 2,4,6-Tribromophenol	15.954	330	51	0.011	ng	0.00
15) 2-Fluorobiphenyl	13.095	172	8244	0.269	ng	0.00
27) Fluoranthene-d10	19.236	212	15404	0.345	ng	0.00
31) Terphenyl-d14	19.830	244	9169	0.412	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.297	149	2529	0.090	ng	# 99

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

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