

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050622\
 Data File : BN019700.D
 Acq On : 06 May 2022 16:08
 Operator : CG/JU
 Sample : N2735-01DL2 100X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 38072-120920DL2

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/09/2022
 Supervised By : mohammad ahmed 05/09/2022

Quant Time: May 07 00:52:02 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN050422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 15:20:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.862	152	5285	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	15607	0.400	ng	0.00
13) Acenaphthene-d10	14.474	164	8498	0.400	ng	-0.01
19) Phenanthrene-d10	17.215	188	16874	0.400	ng	0.00
29) Chrysene-d12	21.404	240	13580	0.400	ng	# 0.00
35) Perylene-d12	23.749	264	10365	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.449	112	52	0.004	ng	0.00
5) Phenol-d6	7.031	99	69	0.005	ng	0.00
8) Nitrobenzene-d5	9.003	82	66	0.005	ng	0.00
11) 2-Methylnaphthalene-d10	12.238	152	83	0.003	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	10	0.003	ng	0.00
15) 2-Fluorobiphenyl	13.116	172	157	0.004	ng	0.01
27) Fluoranthene-d10	19.248	212	156	0.003	ng	0.00
31) Terphenyl-d14	19.847	244	151	0.005	ng	0.00
Target Compounds						Qvalue
37) Benzo(b)fluoranthene	23.039	252	59154m	1.381	ng	

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

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