

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032822\
 Data File : BN019195.D
 Acq On : 28 Mar 2022 14:32
 Operator : CG/JU
 Sample : N1870-05DL 50X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-BN-WPDL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/29/2022
 Supervised By : Jagrut Upadhyay 03/29/2022

Quant Time: Mar 28 15:09:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN032522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 19:03:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	3953	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	10380	0.400	ng	# 0.00
13) Acenaphthene-d10	14.484	164	6239	0.400	ng	0.00
19) Phenanthrene-d10	17.225	188	12902	0.400	ng	# 0.00
29) Chrysene-d12	21.413	240	10130	0.400	ng	# 0.00
35) Perylene-d12	23.790	264	8049	0.400	ng	# 0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.392	112	26813	2.871	ng	0.00
5) Phenol-d6	7.002	99	28194	2.637	ng	0.00
8) Nitrobenzene-d5	8.993	82	13953	1.734	ng	0.00
11) 2-Methylnaphthalene-d10	12.269	152	5	0.000	ng	0.03
14) 2,4,6-Tribromophenol	15.974	330	7779	2.565	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	48010	1.831	ng	0.00
27) Fluoranthene-d10	19.261	212	15	0.000	ng	0.00
31) Terphenyl-d14	19.855	244	43431	1.920	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.550	42	10336	1.757	ng	# 97
12) 2-Methylnaphthalene	12.310	142	50498	2.629	ng	99

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

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