

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032822\
 Data File : BN019194.D
 Acq On : 28 Mar 2022 13:29
 Operator : CG/JU
 Sample : N1870-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-BN-WP

Manual Integrations
 APPROVED

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

Quant Time: Mar 28 14:05:05 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	3714	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	9363	0.400	ng	# 0.00
13) Acenaphthene-d10	14.485	164	5194	0.400	ng	0.00
19) Phenanthrene-d10	17.226	188	10628	0.400	ng	0.00
29) Chrysene-d12	21.422	240	8138	0.400	ng	0.00
35) Perylene-d12	23.782	264	7858	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.392	112	1360462	155.071	ng	0.00
5) Phenol-d6	7.002	99	1757670	174.962	ng	0.00
8) Nitrobenzene-d5	8.993	82	935007	128.852	ng	0.00
11) 2-Methylnaphthalene-d10	12.201	152	2	0.000	ng	-0.04
14) 2,4,6-Tribromophenol	15.974	330	479518	189.930	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	2030612	93.038	ng	0.00
27) Fluoranthene-d10	19.261	212	24	0.001	ng	0.00
31) Terphenyl-d14	19.856	244	1648893	90.756	ng	0.00
Target Compounds						
3) n-Nitrosodimethylamine	3.543	42	517131	93.578	ng	# 99
10) Hexachlorobutadiene	10.989	225	1554	0.263	ng	# 100
12) 2-Methylnaphthalene	12.310	142	2013998	116.220	ng	99

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

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