

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050622\
 Data File : BN019698.D
 Acq On : 06 May 2022 14:01
 Operator : CG/JU
 Sample : N2735-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 38072-120920

Manual Integrations
 APPROVED

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

Quant Time: May 07 00:51:52 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	5896	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	18429	0.400	ng	0.00
13) Acenaphthene-d10	14.485	164	9774	0.400	ng	0.00
19) Phenanthrene-d10	17.215	188	17957	0.400	ng	0.00
29) Chrysene-d12	21.404	240	9695	0.400	ng	0.00
35) Perylene-d12	23.750	264	8629	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.442	112	5237	0.381	ng	0.00
5) Phenol-d6	7.024	99	6269	0.372	ng	0.00
8) Nitrobenzene-d5	9.003	82	5420m	0.363	ng	0.00
11) 2-Methylnaphthalene-d10	12.238	152	12006	0.400	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	964	0.267	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	16429	0.390	ng	0.00
27) Fluoranthene-d10	19.249	212	18118	0.364	ng	0.00
31) Terphenyl-d14	19.847	244	11392	0.502	ng	0.00
Target Compounds						Qvalue
37) Benzo(b)fluoranthene	23.048	252	4602603	129.072	ng	# 89

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

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