

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092221\
 Data File : BN016409.D
 Acq On : 22 Sep 2021 12:59
 Operator : CG/JU
 Sample : M3835-03DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SW-CB-04-(1.0)DL

Manual Integrations
 APPROVED

mohammad
 9/23/2021 1:21:52 PM

Quant Time: Sep 22 13:37:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 13:21:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	25208	0.400	ng	0.00
7) Naphthalene-d8	10.444	136	110254	0.400	ng	0.00
13) Acenaphthene-d10	14.294	164	60158	0.400	ng	-0.01
19) Phenanthrene-d10	17.054	188	119529	0.400	ng	# 0.00
29) Chrysene-d12	21.259	240	103824	0.400	ng	# 0.00
35) Perylene-d12	23.518	264	94591	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.291	112	678	0.011	ng	0.00
5) Phenol-d6	6.849	99	559	0.007	ng	0.00
8) Nitrobenzene-d5	8.809	82	1521m	0.023	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	3786	0.022	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	476	0.027	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	5453	0.023	ng	0.00
27) Fluoranthene-d10	19.092	212	8599	0.025	ng	0.00
31) Terphenyl-d14	19.703	244	6996	0.029	ng	0.00
Target Compounds						
9) Naphthalene	10.495	128	302672	1.019	ng	Qvalue 100

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

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