

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082520\
 Data File : BN011639.D
 Acq On : 26 Aug 2020 05:51
 Operator : CG/JU
 Sample : L3737-04
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 KY038PS038-200818

Quant Time: Aug 26 09:08:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 25 03:31:11 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	2251	0.40	ng	0.00
7) Naphthalene-d8	10.54	136	8915	0.40	ng	0.00
13) Acenaphthene-d10	14.38	164	5556	0.40	ng	-0.01
19) Phenanthrene-d10	17.14	188	12482	0.40	ng	0.00
29) Chrysene-d12	21.32	240	11413	0.40	ng	-0.01
36) Perylene-d12	23.58	264	12116	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.35	112	1119	0.16	ng	0.00
5) Phenol-d6	6.91	99	871	0.09	ng	0.00
8) Nitrobenzene-d5	8.89	82	4213	0.55	ng	-0.01
11) 2-Methylnaphthalene-d10	12.13	152	4664	0.28	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	1086	0.36	ng	0.00
15) 2-Fluorobiphenyl	13.01	172	11910	0.55	ng	-0.01
27) Fluoranthene-d10	19.17	212	12600	0.31	ng	0.00
31) Terphenyl-d14	19.77	244	22370	0.77	ng	0.00
Target Compounds						
30) Pyrene	19.56	202	878	0.022	ng	# 94
32) Benzo(a)anthracene	21.31	228	874	0.021	ng	# 85
33) Chrysene	21.35	228	842	0.021	ng	# 86
37) Benzo(b)fluoranthene	22.90	252	888	0.020	ng	# 27

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

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