

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031120\
 Data File : BN010196.D
 Acq On : 11 Mar 2020 17:35
 Operator : CG/JU
 Sample : L1842-11DL 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE132D6-20200304DL

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:45:58 PM

Quant Time: Mar 12 07:36:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 11 17:37:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	763	0.40	ng	0.04
7) Naphthalene-d8	10.19	136	2786m	0.40	ng	0.08
13) Acenaphthene-d10	14.02	164	1733	0.40	ng	0.02
19) Phenanthrene-d10	16.80	188	3398	0.40	ng	0.03
27) Chrysene-d12	21.00	240	3600	0.40	ng	0.01
34) Perylene-d12	23.10	264	3803	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.11	112	105m	0.06	ng	0.07
5) Phenol-d6	0.00	99	0d	0.00	ng	
8) Nitrobenzene-d5	8.73	82	280m	0.12	ng	0.18
11) 2-Methylnaphthalene-d10	11.81	152	510	0.09	ng	0.08
14) 2,4,6-Tribromophenol	15.58	330	106	0.16	ng	0.04
15) 2-Fluorobiphenyl	12.68	172	921m	0.14	ng	0.04
25) Fluoranthene-d10	18.84	212	1607	0.14	ng	0.02
29) Terphenyl-d14	19.44	244	1389	0.16	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.04	88	2265m	2.898	ng	
23) Phenanthrene	16.85	178	316	0.031	ng	98

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

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