

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031020\
 Data File : BN010167.D
 Acq On : 10 Mar 2020 19:28
 Operator : CG/JU
 Sample : L1842-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP01-20200304

Manual Integrations
 APPROVED

mohammad
 3/11/2020 5:07:53 PM

Quant Time: Mar 11 11:56:48 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 Feb 19 15:57:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	722	0.40	ng	-0.05
7) Naphthalene-d8	10.14	136	2446	0.40	ng	-0.04
13) Acenaphthene-d10	14.01	164	1545	0.40	ng	-0.05
19) Phenanthrene-d10	16.79	188	3115	0.40	ng	-0.03
27) Chrysene-d12	21.00	240	3459	0.40	ng	-0.04
34) Perylene-d12	23.09	264	3780	0.40	ng	-0.05
System Monitoring Compounds						
4) 2-Fluorophenol	5.04	112	348	0.22	ng	-0.04
5) Phenol-d6	6.63	99	223	0.11	ng	-0.02
8) Nitrobenzene-d5	8.59	82	795	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	11.74	152	1631	0.34	ng	-0.04
14) 2,4,6-Tribromophenol	15.56	330	304	0.51	ng	-0.03
15) 2-Fluorobiphenyl	12.64	172	2570	0.44	ng	-0.05
25) Fluoranthene-d10	18.82	212	4000	0.37	ng	-0.04
29) Terphenyl-d14	19.42	244	4360	0.53	ng	-0.05
Target Compounds						
2) 1,4-Dioxane	3.03	88	3193m	4.318	ng	Qvalue
6) bis(2-Chloroethyl)ether	6.86	93	160	0.084	ng	# 82

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

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