

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF093019\
 Data File : BF117266.D
 Acq On : 30 Sep 2019 17:49
 Operator : JU
 Sample : K5019-02DL 20X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW30D-GWDL

Manual Integrations
 APPROVED

mohammad
 10/2/2019 8:32:27 AM

Quant Time: Oct 01 02:05:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	24997	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	113017	20.00	ng	-0.01
39) Acenaphthene-d10	9.85	164	58414	20.00	ng	-0.01
64) Phenanthrene-d10	11.33	188	96072	20.00	ng	0.00
76) Chrysene-d12	13.97	240	78282	20.00	ng	0.00
87) Perylene-d12	15.43	264	75289	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	3327	2.08	ng	0.02
7) Phenol-d6	6.47	99	3000	1.45	ng	0.02
23) Nitrobenzene-d5	7.39	82	2543	1.18	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	791	1.62	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	5855	1.57	ng	0.00
79) Terphenyl-d14	12.91	244	8392	2.00	ng	0.00
Target Compounds						
31) Naphthalene	8.12	128	896940	165.018	ng	99
37) 2-Methylnaphthalene	8.80	142	37759	10.316	ng	96
38) 1-Methylnaphthalene	8.90	142	36949m	10.778	ng	
52) Acenaphthene	9.87	154	15446	4.810	ng	97
55) Dibenzofuran	10.06	168	15727	3.319	ng	95

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

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