

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061219\
 Data File : BN006281.D
 Acq On : 12 Jun 2019 15:34
 Operator : JU/SJ
 Sample : K3231-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8L3

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:33:49 AM

Quant Time: Jun 12 17:29:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4389	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	17256	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9011	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	17385m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	11255m	0.40	ng/ul	0.04
20) Perylene-d12	23.57	264	15635	0.40	ng/ul	0.08
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	3756	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	7862m	0.17	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.72	266	395m	0.089	ng/ul	
12) Phenanthrene	17.09	178	1200m	0.022	ng/ul	
15) Fluoranthene	19.13	202	2451m	0.040	ng/ul	
17) Pyrene	19.49	202	2242	0.038	ng/ul#	78
18) Benzo(a)anthracene	21.29	228	1387	0.028	ng/ul#	77
19) Chrysene	21.34	228	1483	0.032	ng/ul#	83
21) Benzo(b)fluoranthene	22.90	252	2759m	0.042	ng/ul	

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

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