

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061219\
 Data File : BN006290.D
 Acq On : 12 Jun 2019 20:46
 Operator : JU/SJ
 Sample : K3231-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8L8

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:34:10 AM

Quant Time: Jun 13 02:49:25 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 : Wed Jun 12 17:24:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	5394	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	21761	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11368	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	21675m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	13198m	0.40	ng/ul	-0.02
20) Perylene-d12	23.50	264	12955m	0.40	ng/ul	-0.05
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5320	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	9171m	0.16	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.71	266	276m	0.050	ng/ul	
12) Phenanthrene	17.09	178	1818m	0.026	ng/ul	
15) Fluoranthene	19.12	202	4742m	0.062	ng/ul	
17) Pyrene	19.49	202	3846m	0.055	ng/ul	
18) Benzo(a)anthracene	21.25	228	1543m	0.026	ng/ul	
19) Chrysene	21.30	228	2326m	0.042	ng/ul	
21) Benzo(b)fluoranthene	22.83	252	3814m	0.071	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	1685m	0.032	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	1653	0.030	ng/ul#	95
26) Benzo(g,h,i)perylene	26.42	276	1562	0.034	ng/ul#	50

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

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