

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004049.D
 Acq On : 17 Dec 2018 23:07
 Operator : JU/SJ
 Sample : J6236-19
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F19

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:29 PM

Quant Time: Dec 18 06:21:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 18 04:29:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1445	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7731	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5129	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13211m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	15022m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12994m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3878	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	13465m	0.35	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.20	152	846	0.037	ng/ul#	74
12) Phenanthrene	17.26	178	1201m	0.028	ng/ul	
15) Fluoranthene	19.27	202	2128m	0.042	ng/ul	
17) Pyrene	19.64	202	2336	0.043	ng/ul#	86
18) Benzo(a)anthracene	21.39	228	1903	0.040	ng/ul#	82
19) Chrysene	21.44	228	2204	0.041	ng/ul#	83
21) Benzo(b)fluoranthene	23.02	252	3266m	0.061	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1123m	0.021	ng/ul	
23) Benzo(a)pyrene	23.62	252	1677	0.036	ng/ul#	41
24) Indeno(1,2,3-cd)pyrene	26.09	276	1405	0.026	ng/ul#	99
26) Benzo(g,h,i)perylene	26.82	276	1080	0.023	ng/ul#	48

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

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