

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004096.D
 Acq On : 19 Dec 2018 02:39
 Operator : JU/SJ
 Sample : J6271-17
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E32

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:53 PM

Quant Time: Dec 19 04:54:57 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 : Wed Dec 19 04:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2086	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10173	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6007	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	14900m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	15507m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13714m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3139	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	9381m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	877	0.032	ng/ul#	80
12) Phenanthrene	17.25	178	1761m	0.036	ng/ul	
15) Fluoranthene	19.27	202	5059m	0.088	ng/ul	
17) Pyrene	19.64	202	5286m	0.095	ng/ul	
18) Benzo(a)anthracene	21.38	228	3943	0.081	ng/ul#	90
19) Chrysene	21.43	228	5389	0.096	ng/ul#	92
21) Benzo(b)fluoranthene	23.02	252	7164m	0.126	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	2942m	0.053	ng/ul	
23) Benzo(a)pyrene	23.62	252	3800	0.077	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	26.08	276	3386	0.060	ng/ul#	94
26) Benzo(g,h,i)perylene	26.82	276	2972	0.061	ng/ul#	72

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

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