

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN003997.D
 Acq On : 16 Dec 2018 16:22
 Operator : JU/SJ
 Sample : J6229-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F30

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:27:40 PM

Quant Time: Dec 17 05:04:30 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 : Mon Dec 17 04:06:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2047	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10114	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6269	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	15121m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	17115m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	15417m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5202	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	15874m	0.36	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.29	142	416	0.021	ng/ul	97
12) Phenanthrene	17.26	178	7646	0.154	ng/ul	98
15) Fluoranthene	19.28	202	2964m	0.051	ng/ul	
17) Pyrene	19.64	202	7875m	0.128	ng/ul	
18) Benzo(a)anthracene	21.39	228	3653m	0.068	ng/ul	
19) Chrysene	21.44	228	11456m	0.185	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	4785m	0.075	ng/ul	
23) Benzo(a)pyrene	23.62	252	3451m	0.062	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.09	276	2238m	0.035	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	1533m	0.030	ng/ul	
26) Benzo(g,h,i)perylene	26.82	276	3264m	0.059	ng/ul	

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

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