

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN004011.D
 Acq On : 17 Dec 2018 00:40
 Operator : JU/SJ
 Sample : J6228-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E58

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:28:19 PM

Quant Time: Dec 17 07:01:22 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 06:22:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1990	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9883	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5911	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14715m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16388m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13936m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3731	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	11549m	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1873	0.067	ng/ul#	86
5) 2-Methylnaphthalene	12.29	142	3713	0.192	ng/ul	98
9) Fluorene	15.52	166	821	0.030	ng/ul#	96
12) Phenanthrene	17.26	178	87652m	1.815	ng/ul	
13) Anthracene	17.35	178	5572m	0.140	ng/ul	
15) Fluoranthene	19.27	202	19099m	0.335	ng/ul	
17) Pyrene	19.64	202	85664	1.457	ng/ul	96
18) Benzo(a)anthracene	21.38	228	47060m	0.909	ng/ul	
19) Chrysene	21.43	228	83659m	1.412	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	25713m	0.446	ng/ul	
23) Benzo(a)pyrene	23.62	252	32753m	0.656	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	18138m	0.317	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	14090m	0.306	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	37713m	0.758	ng/ul	

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

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