

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004092.D
 Acq On : 19 Dec 2018 00:19
 Operator : JU/SJ
 Sample : J6271-13
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E28

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:46 PM

Quant Time: Dec 19 04:41:50 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	1748	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8478	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5148	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13111m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14596m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13338m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2846	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	9062m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	38937	1.680	ng/ul	100
9) Fluorene	15.52	166	824	0.035	ng/ul#	94
12) Phenanthrene	17.26	178	9247m	0.215	ng/ul	
13) Anthracene	17.35	178	33535m	0.948	ng/ul	
15) Fluoranthene	19.27	202	174923m	3.442	ng/ul	
17) Pyrene	19.64	202	189508	3.620	ng/ul	98
18) Benzo(a)anthracene	21.38	228	132149m	2.867	ng/ul	
19) Chrysene	21.44	228	130352m	2.471	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	257983m	4.676	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	82728m	1.539	ng/ul	
23) Benzo(a)pyrene	23.62	252	116360m	2.435	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	107931m	1.970	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	26702m	0.606	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	82318m	1.729	ng/ul	

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

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