

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004143.D
 Acq On : 20 Dec 2018 08:30
 Operator : JU/SJ
 Sample : J6299-04
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F23

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:04:17 PM

Quant Time: Dec 20 11:21:05 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 : Thu Dec 20 07:15:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1103	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	5886	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	3659	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	9408	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	10083	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	9809	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2032	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	6393	0.23	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	6311	0.383	ng/ul	98
12) Phenanthrene	17.25	178	1735m	0.056	ng/ul	
13) Anthracene	17.35	178	6051	0.238	ng/ul	97
15) Fluoranthene	19.27	202	29548	0.810	ng/ul	97
17) Pyrene	19.63	202	30733	0.850	ng/ul	95
18) Benzo(a)anthracene	21.38	228	19959	0.627	ng/ul	97
19) Chrysene	21.43	228	20218	0.555	ng/ul	98
21) Benzo(b)fluoranthene	23.02	252	34800m	0.858	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	12730m	0.322	ng/ul	
23) Benzo(a)pyrene	23.62	252	15078	0.429	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.08	276	13198	0.328	ng/ul	95
25) Dibenzo(a,h)anthracene	26.08	278	3612	0.111	ng/ul#	76
26) Benzo(g,h,i)perylene	26.81	276	10185	0.291	ng/ul#	93

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

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