

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003979.D
 Acq On : 16 Dec 2018 03:32
 Operator : JU/SJ
 Sample : J6171-13DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y17DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:11 PM

Quant Time: Dec 17 00:14:07 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 : Sat Dec 15 23:04:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1791	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9299	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5706	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14442	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	18865	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	17256	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	701	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	2270	0.05	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.68	128	1115	0.043	ng/ul#	80
5) 2-Methylnaphthalene	12.29	142	596	0.033	ng/ul	99
12) Phenanthrene	17.26	178	3345	0.071	ng/ul#	94
17) Pyrene	19.63	202	5304	0.078	ng/ul#	92
18) Benzo(a)anthracene	21.38	228	1934	0.032	ng/ul#	56
19) Chrysene	21.43	228	4315	0.063	ng/ul#	80
21) Benzo(b)fluoranthene	23.01	252	1709m	0.024	ng/ul	
23) Benzo(a)pyrene	23.61	252	1664	0.027	ng/ul#	4
26) Benzo(g,h,i)perylene	26.79	276	3223	0.052	ng/ul#	63

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

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