

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004081.D
 Acq On : 18 Dec 2018 17:53
 Operator : JU/SJ
 Sample : J6271-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F11

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:04 PM

Quant Time: Dec 19 03:52:06 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 01:54:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1459	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7325	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4917	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12899m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13952m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12757m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2455	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	8366m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.29	142	3960	0.276	ng/ul	100
8) Acenaphthene	14.53	153	1743	0.090	ng/ul#	88
9) Fluorene	15.52	166	2614	0.116	ng/ul#	91
12) Phenanthrene	17.26	178	22779	0.538	ng/ul	96
13) Anthracene	17.35	178	2685m	0.077	ng/ul	
15) Fluoranthene	19.27	202	2983m	0.060	ng/ul	
17) Pyrene	19.64	202	21067m	0.421	ng/ul	
18) Benzo(a)anthracene	21.39	228	3210m	0.073	ng/ul	
19) Chrysene	21.44	228	7785m	0.154	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	1800m	0.034	ng/ul	
23) Benzo(a)pyrene	23.62	252	1871	0.041	ng/ul#	8
26) Benzo(g,h,i)perylene	26.82	276	1599	0.035	ng/ul#	46

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

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