

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004057.D
 Acq On : 18 Dec 2018 03:49
 Operator : JU/SJ
 Sample : J6270-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E04

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:51 PM

Quant Time: Dec 18 07:25:18 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 : Tue Dec 18 06:35:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1642	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8168	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5037	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13059m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14687m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12855m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3266	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	10910m	0.29	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	1712	0.075	ng/ul#	89
12) Phenanthrene	17.26	178	2525m	0.059	ng/ul	
13) Anthracene	17.35	178	1769m	0.050	ng/ul	
15) Fluoranthene	19.28	202	8716m	0.172	ng/ul	
17) Pyrene	19.64	202	9590m	0.182	ng/ul	
18) Benzo(a)anthracene	21.38	228	6630m	0.143	ng/ul	
19) Chrysene	21.44	228	8426m	0.159	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	12691m	0.239	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	3823m	0.074	ng/ul	
23) Benzo(a)pyrene	23.62	252	6010m	0.131	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.09	276	5014m	0.095	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	1458m	0.034	ng/ul	
26) Benzo(g,h,i)perylene	26.82	276	3899m	0.085	ng/ul	

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

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