

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004053.D
 Acq On : 18 Dec 2018 01:28
 Operator : JU/SJ
 Sample : J6270-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E00

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:39 PM

Quant Time: Dec 18 06:56:28 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	7.83	152	1581	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7683	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4675	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12231m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14543m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13159m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3344	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	11656m	0.33	ng/ul	0.00
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.29	142	327	0.022	ng/ul	98
7) Acenaphthylene	14.19	152	17192	0.817	ng/ul	99
9) Fluorene	15.52	166	900	0.042	ng/ul#	95
12) Phenanthrene	17.25	178	6575m	0.164	ng/ul	
13) Anthracene	17.35	178	16271m	0.493	ng/ul	
15) Fluoranthene	19.27	202	64557m	1.362	ng/ul	
17) Pyrene	19.64	202	72931	1.398	ng/ul	98
18) Benzo(a)anthracene	21.38	228	50172m	1.092	ng/ul	
19) Chrysene	21.43	228	60228m	1.146	ng/ul	
21) Benzo(b)fluoranthene	23.01	252	118003m	2.168	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	38574m	0.727	ng/ul	
23) Benzo(a)pyrene	23.62	252	51182m	1.086	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	55402m	1.025	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	14242m	0.327	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	64004m	1.362	ng/ul	

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

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