

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004061.D
 Acq On : 18 Dec 2018 06:09
 Operator : JU/SJ
 Sample : J6270-09
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E08

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:57:03 PM

Quant Time: Dec 18 08:14:13 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	1659	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8452	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5363	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14144m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16442m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	15581m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3109	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	10560m	0.26	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	8338	0.345	ng/ul	98
12) Phenanthrene	17.26	178	4088m	0.088	ng/ul	
13) Anthracene	17.35	178	6794m	0.178	ng/ul	
15) Fluoranthene	19.28	202	38418m	0.701	ng/ul	
17) Pyrene	19.64	202	43291	0.734	ng/ul	95
18) Benzo(a)anthracene	21.39	228	26823m	0.517	ng/ul	
19) Chrysene	21.44	228	31317m	0.527	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	56686m	0.879	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	18003m	0.287	ng/ul	
23) Benzo(a)pyrene	23.62	252	26572	0.476	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.08	276	24132	0.377	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.08	278	5877	0.114	ng/ul#	58
26) Benzo(g,h,i)perylene	26.82	276	21553	0.387	ng/ul#	91

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

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