

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004062.D
 Acq On : 18 Dec 2018 06:44
 Operator : JU/SJ
 Sample : J6270-10
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E09

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:57:05 PM

Quant Time: Dec 18 08:17:48 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	1752	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8882	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5523	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14276m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16059m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	14859m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3104	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	10471m	0.25	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	3844	0.155	ng/ul	95
12) Phenanthrene	17.26	178	2918	0.062	ng/ul	93
13) Anthracene	17.35	178	3220m	0.084	ng/ul	
15) Fluoranthene	19.28	202	20138m	0.364	ng/ul	
17) Pyrene	19.64	202	21235m	0.369	ng/ul	
18) Benzo(a)anthracene	21.39	228	14015	0.276	ng/ul	94
19) Chrysene	21.44	228	15291m	0.263	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	27019m	0.440	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	9052m	0.151	ng/ul	
23) Benzo(a)pyrene	23.62	252	12382	0.233	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.08	276	10926	0.179	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.08	278	2671	0.054	ng/ul#	32
26) Benzo(g,h,i)perylene	26.82	276	9533	0.180	ng/ul#	84

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

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