

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004097.D
 Acq On : 19 Dec 2018 03:14
 Operator : JU/SJ
 Sample : J6271-18
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E33

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:54 PM

Quant Time: Dec 19 04:57:36 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 04:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1981	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9636	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5789	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	14459m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	15092m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13637m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2840	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	8773m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	544	0.021	ng/ul#	67
12) Phenanthrene	17.25	178	1718m	0.036	ng/ul	
15) Fluoranthene	19.27	202	2696m	0.048	ng/ul	
17) Pyrene	19.64	202	2942	0.054	ng/ul#	90
18) Benzo(a)anthracene	21.38	228	1937	0.041	ng/ul#	85
19) Chrysene	21.43	228	2637	0.048	ng/ul#	88
21) Benzo(b)fluoranthene	23.02	252	3313m	0.059	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1125m	0.020	ng/ul	
23) Benzo(a)pyrene	23.62	252	1814	0.037	ng/ul#	46
24) Indeno(1,2,3-cd)pyrene	26.08	276	1596	0.028	ng/ul#	95
26) Benzo(g,h,i)perylene	26.81	276	1419	0.029	ng/ul#	54

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

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