

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004114.D
 Acq On : 19 Dec 2018 15:32
 Operator : JU/SJ
 Sample : J6297-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F26

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:03:05 PM

Quant Time: Dec 20 05:49:11 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 : Thu Dec 20 04:42:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1356	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7041	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4497	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11797m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12086m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	10402m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2597	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	8562m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.26	178	1195m	0.031	ng/ul	
15) Fluoranthene	19.27	202	1640m	0.036	ng/ul	
17) Pyrene	19.63	202	2220	0.051	ng/ul#	82
18) Benzo(a)anthracene	21.38	228	1105	0.029	ng/ul#	78
19) Chrysene	21.43	228	1504	0.034	ng/ul#	85
21) Benzo(b)fluoranthene	23.01	252	2002m	0.047	ng/ul	
23) Benzo(a)pyrene	23.62	252	994	0.027	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	26.08	276	973	0.023	ng/ul#	96
26) Benzo(g,h,i)perylene	26.81	276	1132	0.030	ng/ul#	54

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

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