

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN004008.D
 Acq On : 16 Dec 2018 22:53
 Operator : JU/SJ
 Sample : J6228-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E86

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:28:15 PM

Quant Time: Dec 17 06:45:07 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 : Mon Dec 17 06:22:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1811	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9234	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5645	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14290m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16328m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	14151m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2513	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	8033m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.26	178	3409	0.073	ng/ul	95
15) Fluoranthene	19.28	202	4199m	0.076	ng/ul	
17) Pyrene	19.64	202	5507	0.094	ng/ul#	91
18) Benzo(a)anthracene	21.38	228	3304m	0.064	ng/ul	
19) Chrysene	21.44	228	4624m	0.078	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	3604m	0.062	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1457m	0.026	ng/ul	
23) Benzo(a)pyrene	23.62	252	2580m	0.051	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	1870m	0.032	ng/ul	
26) Benzo(g,h,i)perylene	26.82	276	2255m	0.045	ng/ul	

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

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