

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN004019.D
 Acq On : 17 Dec 2018 05:24
 Operator : JU/SJ
 Sample : J6236-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E98

Manual Integrations
 APPROVED

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

Quant Time: Dec 17 07:54:05 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	2130	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10691	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6463	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	16299m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	17216m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	14999m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3093	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	9500m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.29	142	722	0.034	ng/ul	99
12) Phenanthrene	17.26	178	11579	0.216	ng/ul	99
15) Fluoranthene	19.27	202	8943m	0.142	ng/ul	
17) Pyrene	19.64	202	13345m	0.216	ng/ul	
18) Benzo(a)anthracene	21.38	228	6669m	0.123	ng/ul	
19) Chrysene	21.44	228	15145	0.243	ng/ul	96
21) Benzo(b)fluoranthene	23.01	252	8508m	0.137	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	2428m	0.040	ng/ul	
23) Benzo(a)pyrene	23.62	252	5659	0.105	ng/ul#	69
24) Indeno(1,2,3-cd)pyrene	26.08	276	3962	0.064	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.08	278	1878	0.038	ng/ul#	20
26) Benzo(g,h,i)perylene	26.81	276	5220	0.097	ng/ul#	79

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

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