

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN112018\
 Data File : BN003627.D
 Acq On : 20 Nov 2018 12:21
 Operator : MJ/SJ
 Sample : J6001-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40T2

Quant Time: Nov 20 13:18:55 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 Nov 20 03:27:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5040	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20019	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	10944	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	26697	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	29110	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	26357	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	11436	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	30072	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.69	128	13459	0.239	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	10120	0.258	ng/ul	100
7) Acenaphthylene	14.20	152	38212	0.775	ng/ul	100
8) Acenaphthene	14.54	153	13560	0.314	ng/ul	99
9) Fluorene	15.53	166	17093	0.340	ng/ul	99
11) Pentachlorophenol	16.87	266	996	0.207	ng/ul	99
12) Phenanthrene	17.26	178	59893	0.684	ng/ul	100
13) Anthracene	17.36	178	16663	0.231	ng/ul	100
15) Fluoranthene	19.28	202	42072	0.407	ng/ul	100
17) Pyrene	19.64	202	45860	0.439	ng/ul	100
18) Benzo(a)anthracene	21.38	228	64833	0.705	ng/ul	100
19) Chrysene	21.43	228	39660	0.377	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	35863	0.329	ng/ul	99
23) Benzo(a)pyrene	23.61	252	44231	0.468	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	37092	0.343	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.08	278	31670	0.363	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	37610	0.400	ng/ul	99

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

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