

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112018\
 Data File : BN003626.D
 Acq On : 20 Nov 2018 11:05
 Operator : MJ/SJ
 Sample : J6002-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2E83

Quant Time: Nov 20 11:50:59 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	5002	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20724	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11070	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	28028	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	30770	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	30020	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	8249	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	32334	0.39	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.69	128	49523	0.851	ng/ul 100
7) Acenaphthylene	14.20	152	22362	0.449	ng/ul 100
8) Acenaphthene	14.54	153	21271	0.487	ng/ul 99
9) Fluorene	15.53	166	20653	0.406	ng/ul 99
11) Pentachlorophenol	16.86	266	5437	1.077	ng/ul 98
12) Phenanthrene	17.26	178	47477	0.516	ng/ul 100
13) Anthracene	17.35	178	42648	0.564	ng/ul 100
15) Fluoranthene	19.27	202	62998	0.580	ng/ul 95
17) Pyrene	19.64	202	83628	0.758	ng/ul 99
18) Benzo(a)anthracene	21.38	228	62781	0.646	ng/ul 99
19) Chrysene	21.38	228	62781	0.564	ng/ul 96
21) Benzo(b)fluoranthene	23.01	252	65793	0.530	ng/ul 99
22) Benzo(k)fluoranthene	23.01	252	65793	0.544	ng/ul 99
23) Benzo(a)pyrene	23.61	252	4719	0.044	ng/ul# 63
24) Indeno(1,2,3-cd)pyrene	26.06	276	89995	0.730	ng/ul 99
25) Dibenzo(a,h)anthracene	26.07	278	69719	0.702	ng/ul 99
26) Benzo(g,h,i)perylene	26.78	276	76088	0.710	ng/ul 99

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

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