

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004262.D
 Acq On : 28 Dec 2018 15:01
 Operator : JU/SJ
 Sample : SST0.855
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.855

Quant Time: Dec 28 15:59:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 15:21:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2143	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	9009	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.46	164	5311	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	11790	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	14428	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	12461	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	11714	0.99	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	28313	1.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.67	128	20681	0.717	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	14885	0.813	ng/ul	100
7) Acenaphthylene	14.19	152	21610	0.818	ng/ul	99
8) Acenaphthene	14.53	153	16879	0.662	ng/ul	100
9) Fluorene	15.52	166	20248	0.731	ng/ul	100
11) Pentachlorophenol	16.88	266	1347	0.568	ng/ul	97
12) Phenanthrene	17.26	178	31104	0.646	ng/ul	99
13) Anthracene	17.35	178	29184	0.816	ng/ul	99
15) Fluoranthene	19.28	202	37616	0.806	ng/ul	99
17) Pyrene	19.64	202	38776	0.457	ng/ul	99
18) Benzo(a)anthracene	21.39	228	38633	0.682	ng/ul	99
19) Chrysene	21.44	228	40074	0.520	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	38684	0.516	ng/ul	98
22) Benzo(k)fluoranthene	23.07	252	39221	0.555	ng/ul	98
23) Benzo(a)pyrene	23.63	252	35558	0.606	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.11	276	37284	0.563	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.11	278	30598	0.640	ng/ul	97
26) Benzo(g,h,i)perylene	26.83	276	30534	0.468	ng/ul	98

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

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