

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005147.D
 Acq On : 18 Apr 2019 10:14
 Operator : JU/SJ
 Sample : SST0.855
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.855

Quant Time: Apr 18 11:19:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 11:12:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	3199	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	11790	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	6653	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	15417	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	12822	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	11229	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	15414	0.83	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	35645	0.74	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.42	128	26283	0.776	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	18798	0.777	ng/ul	100
7) Acenaphthylene	13.95	152	26794	0.881	ng/ul	99
8) Acenaphthene	14.30	153	20171	0.764	ng/ul	99
9) Fluorene	15.30	166	23477	0.744	ng/ul	99
11) Pentachlorophenol	16.64	266	2867	0.873	ng/ul	97
12) Phenanthrene	17.03	178	38631	0.771	ng/ul	100
13) Anthracene	17.12	178	35719	0.846	ng/ul	99
15) Fluoranthene	19.06	202	44476	0.724	ng/ul	99
17) Pyrene	19.42	202	43755	1.060	ng/ul	99
18) Benzo(a)anthracene	21.19	228	39214	0.948	ng/ul	99
19) Chrysene	21.24	228	39701	0.827	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	35609	0.793	ng/ul	97
22) Benzo(k)fluoranthene	22.78	252	35230	0.798	ng/ul	98
23) Benzo(a)pyrene	23.29	252	31689	0.789	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.56	276	34262	0.649	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.58	278	28292	0.662	ng/ul	98
26) Benzo(g,h,i)perylene	26.21	276	29027	0.606	ng/ul	100

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

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