

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004260.D
 Acq On : 28 Dec 2018 13:51
 Operator : JU/SJ
 Sample : SST0.253
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.253

Manual Integrations
 APPROVED

Sohil
 1/2/2019 3:39:20 PM

Quant Time: Dec 28 15:26:57 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	2277	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	9919	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5590	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	13274	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15080	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	13292	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3011	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	7414	0.24	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	5501	0.173	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	3800	0.189	ng/ul	97
7) Acenaphthylene	14.20	152	5524	0.199	ng/ul	98
8) Acenaphthene	14.54	153	4408	0.164	ng/ul	99
9) Fluorene	15.53	166	4946	0.170	ng/ul	99
11) Pentachlorophenol	16.91	266	290	0.109	ng/ul	91
12) Phenanthrene	17.27	178	8214	0.152	ng/ul	98
13) Anthracene	17.36	178	7488	0.186	ng/ul	98
15) Fluoranthene	19.28	202	9958	0.189	ng/ul	99
17) Pyrene	19.65	202	10375	0.117	ng/ul	100
18) Benzo(a)anthracene	21.39	228	10045	0.170	ng/ul	99
19) Chrysene	21.44	228	11106	0.138	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	9883	0.124	ng/ul	95
22) Benzo(k)fluoranthene	23.07	252	10449m	0.139	ng/ul	
23) Benzo(a)pyrene	23.64	252	9505	0.152	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	26.11	276	9799	0.139	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.11	278	8061	0.158	ng/ul	94
26) Benzo(g,h,i)perylene	26.84	276	8108	0.117	ng/ul	96

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

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