

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004263.D
 Acq On : 28 Dec 2018 15:36
 Operator : JU/SJ
 Sample : SSTD1.656
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.656

Quant Time: Dec 28 16:19:30 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	2062	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	8893	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.46	164	4996	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	12003	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13426	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	12256	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	22497	1.92	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	60497	2.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.67	128	40359	1.417	ng/ul	98
5) 2-Methylnaphthalene	12.28	142	29101	1.611	ng/ul	100
7) Acenaphthylene	14.19	152	42900	1.726	ng/ul	99
8) Acenaphthene	14.52	153	33917	1.414	ng/ul	99
9) Fluorene	15.51	166	40271	1.546	ng/ul	99
11) Pentachlorophenol	16.88	266	3015	1.249	ng/ul	99
12) Phenanthrene	17.25	178	63832	1.303	ng/ul	99
13) Anthracene	17.35	178	58847	1.617	ng/ul	99
15) Fluoranthene	19.27	202	80929	1.703	ng/ul	99
17) Pyrene	19.64	202	79689	1.008	ng/ul	99
18) Benzo(a)anthracene	21.39	228	75149	1.426	ng/ul	99
19) Chrysene	21.44	228	75695	1.055	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	75342	1.021	ng/ul	97
22) Benzo(k)fluoranthene	23.07	252	76239	1.097	ng/ul	96
23) Benzo(a)pyrene	23.63	252	74124	1.285	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	26.10	276	73800	1.133	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.10	278	60323	1.284	ng/ul	96
26) Benzo(g,h,i)perylene	26.83	276	65687	1.024	ng/ul	97

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

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