

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006258.D
 Acq On : 11 Jun 2019 22:06
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.499

Quant Time: Jun 12 01:12:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	6074	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	24436	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11599	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	20385	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	13456	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	15371	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	14526	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	20612	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	27730	0.411	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	18193	0.387	ng/ul	99
7) Acenaphthylene	14.01	152	25715	0.399	ng/ul	100
8) Acenaphthene	14.36	153	18464	0.413	ng/ul	99
9) Fluorene	15.35	166	18678	0.357	ng/ul	98
11) Pentachlorophenol	16.71	266	1679	0.324	ng/ul	97
12) Phenanthrene	17.09	178	25852	0.395	ng/ul	99
13) Anthracene	17.18	178	22624	0.362	ng/ul	99
15) Fluoranthene	19.12	202	27143	0.376	ng/ul	96
17) Pyrene	19.49	202	28339	0.400	ng/ul	100
18) Benzo(a)anthracene	21.25	228	22598	0.380	ng/ul	99
19) Chrysene	21.30	228	26139	0.468	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	24040	0.375	ng/ul	98
22) Benzo(k)fluoranthene	22.87	252	26152	0.419	ng/ul	97
23) Benzo(a)pyrene	23.40	252	24628	0.413	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.73	276	28590	0.442	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.74	278	22543	0.436	ng/ul	94
26) Benzo(g,h,i)perylene	26.41	276	24544	0.454	ng/ul	96

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

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