

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006161.D
 Acq On : 07 Jun 2019 07:32
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.498

Quant Time: Jun 07 08:11:35 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 : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5213	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	20484	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10391	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	21608	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	18092	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	19680	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	11525	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	23315	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	22021	0.390	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	14511	0.368	ng/ul	98
7) Acenaphthylene	14.01	152	22348	0.387	ng/ul	99
8) Acenaphthene	14.36	153	15727	0.393	ng/ul	100
9) Fluorene	15.36	166	16819	0.359	ng/ul	98
11) Pentachlorophenol	16.72	266	1929	0.351	ng/ul	97
12) Phenanthrene	17.09	178	26056	0.376	ng/ul	100
13) Anthracene	17.19	178	23720	0.358	ng/ul	99
15) Fluoranthene	19.12	202	31219	0.408	ng/ul	96
17) Pyrene	19.49	202	32682	0.343	ng/ul	99
18) Benzo(a)anthracene	21.26	228	27524	0.344	ng/ul	98
19) Chrysene	21.31	228	32099	0.427	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	28891	0.352	ng/ul	98
22) Benzo(k)fluoranthene	22.88	252	31489	0.394	ng/ul	96
23) Benzo(a)pyrene	23.41	252	29853	0.391	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.75	276	34217	0.414	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.76	278	27362	0.414	ng/ul	94
26) Benzo(g,h,i)perylene	26.42	276	29111	0.420	ng/ul	96

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

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