

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009046.D
 Acq On : 20 Dec 2019 01:25
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Dec 20 05:06:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 19 15:54:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	2689	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.28	136	10303	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.16	164	6410	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	14526	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	9956	0.40	ng/ul	0.00
20) Perylene-d12	23.26	264	8428	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	7092	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.95	212	17039	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.33	128	12767	0.409	ng/ul	100
5) 2-Methylnaphthalene	11.96	142	9174	0.389	ng/ul	100
7) Acenaphthylene	13.88	152	13145	0.390	ng/ul	99
8) Acenaphthene	14.23	153	10294	0.390	ng/ul	99
9) Fluorene	15.22	166	12250	0.392	ng/ul	99
11) Pentachlorophenol	16.57	266	1234	0.319	ng/ul	98
12) Phenanthrene	16.95	178	20646	0.409	ng/ul	100
13) Anthracene	17.05	178	18565	0.410	ng/ul	100
15) Fluoranthene	18.98	202	22892	0.396	ng/ul	97
17) Pyrene	19.34	202	22606	0.445	ng/ul	97
18) Benzo(a)anthracene	21.10	228	18704	0.423	ng/ul	99
19) Chrysene	21.15	228	18902	0.425	ng/ul	100
21) Benzo(b)fluoranthene	22.62	252	16107	0.409	ng/ul	98
22) Benzo(k)fluoranthene	22.67	252	17111	0.431	ng/ul	96
23) Benzo(a)pyrene	23.16	252	14912	0.432	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.36	276	15967	0.425	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.37	278	12741	0.426	ng/ul	98
26) Benzo(g,h,i)perylene	25.99	276	13529	0.428	ng/ul	99

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

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