

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101119\
 Data File : BN008067.D
 Acq On : 11 Oct 2019 08:58
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Oct 11 18:59:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 16:11:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2409	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	11506	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	7411	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	17561	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	18075	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	22312	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	7041	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	21616	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	11892	0.363	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	8088	0.360	ng/ul	99
7) Acenaphthylene	13.98	152	12605	0.357	ng/ul	99
8) Acenaphthene	14.33	153	11023	0.402	ng/ul	99
9) Fluorene	15.32	166	11236	0.360	ng/ul	98
11) Pentachlorophenol	16.68	266	1315	0.310	ng/ul	99
12) Phenanthrene	17.06	178	18336	0.357	ng/ul	99
13) Anthracene	17.15	178	15742	0.348	ng/ul	100
15) Fluoranthene	19.08	202	23814	0.351	ng/ul	98
17) Pyrene	19.44	202	27880	0.387	ng/ul	97
18) Benzo(a)anthracene	21.20	228	22828	0.347	ng/ul	99
19) Chrysene	21.25	228	29758	0.369	ng/ul	100
21) Benzo(b)fluoranthene	22.75	252	25063	0.345	ng/ul	98
22) Benzo(k)fluoranthene	22.79	252	30032	0.365	ng/ul	98
23) Benzo(a)pyrene	23.31	252	27095	0.354	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.58	276	32704	0.363	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.60	278	25742	0.362	ng/ul	99
26) Benzo(g,h,i)perylene	26.24	276	29962	0.363	ng/ul	99

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

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