

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN120319\
 Data File : BN008823.D
 Acq On : 03 Dec 2019 09:47
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.424

Quant Time: Dec 03 11:07:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 03 11:01:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	4191	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	17327	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	9917	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	20020	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	16392	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	14496	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	10974	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	23978	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.39	128	20164	0.386	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	14163	0.388	ng/ul	100
7) Acenaphthylene	13.93	152	20568	0.375	ng/ul	100
8) Acenaphthene	14.27	153	14880	0.366	ng/ul	99
9) Fluorene	15.26	166	16686	0.349	ng/ul	99
11) Pentachlorophenol	16.62	266	2108	0.278	ng/ul	100
12) Phenanthrene	17.00	178	27266	0.404	ng/ul	100
13) Anthracene	17.09	178	25963	0.407	ng/ul	99
15) Fluoranthene	19.02	202	31850	0.384	ng/ul	97
17) Pyrene	19.39	202	32507	0.494	ng/ul	98
18) Benzo(a)anthracene	21.15	228	30074	0.456	ng/ul	99
19) Chrysene	21.19	228	28357	0.438	ng/ul	98
21) Benzo(b)fluoranthene	22.68	252	25344	0.424	ng/ul	95
22) Benzo(k)fluoranthene	22.72	252	24578	0.415	ng/ul	96
23) Benzo(a)pyrene	23.23	252	22922	0.409	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.45	276	23239	0.354	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.46	278	18906	0.357	ng/ul	95
26) Benzo(g,h,i)perylene	26.09	276	19239	0.350	ng/ul	98

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

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