

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120219\
 Data File : BN008821.D
 Acq On : 02 Dec 2019 19:05
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.423

Quant Time: Dec 03 00:14:41 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 : Mon Dec 02 12:46:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	5430	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	21897	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	12018	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	23941	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	17618	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	15219	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	13619	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	28258	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.39	128	25553	0.387	ng/ul	99
5) 2-Methylnaphthalene	12.02	142	17445	0.378	ng/ul	98
7) Acenaphthylene	13.93	152	26032	0.392	ng/ul	99
8) Acenaphthene	14.27	153	18020	0.365	ng/ul	99
9) Fluorene	15.27	166	20059	0.346	ng/ul	100
11) Pentachlorophenol	16.62	266	3024	0.334	ng/ul	98
12) Phenanthrene	17.00	178	32246	0.399	ng/ul	99
13) Anthracene	17.09	178	31921	0.418	ng/ul	100
15) Fluoranthene	19.02	202	37469	0.378	ng/ul	100
17) Pyrene	19.39	202	37999	0.538	ng/ul	99
18) Benzo(a)anthracene	21.15	228	32285	0.456	ng/ul	96
19) Chrysene	21.20	228	29812	0.428	ng/ul	96
21) Benzo(b)fluoranthene	22.68	252	26365	0.420	ng/ul	89
22) Benzo(k)fluoranthene	22.72	252	25573	0.411	ng/ul#	89
23) Benzo(a)pyrene	23.23	252	24238	0.411	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.45	276	24924	0.362	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.46	278	20179	0.363	ng/ul#	89
26) Benzo(g,h,i)perylene	26.10	276	21048	0.365	ng/ul	93

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

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