

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN004032.D
 Acq On : 17 Dec 2018 13:04
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.424

Quant Time: Dec 17 14:53:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 17 06:22:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1526	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8009	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4753	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	11825	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13272	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	11893	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5091	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	15495	0.45	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	9018	0.401	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	6668	0.424	ng/ul	100
7) Acenaphthylene	14.19	152	10337	0.483	ng/ul	99
8) Acenaphthene	14.54	153	7702	0.411	ng/ul	99
9) Fluorene	15.52	166	8810	0.403	ng/ul	99
11) Pentachlorophenol	16.87	266	1006	0.472	ng/ul	96
12) Phenanthrene	17.26	178	15341	0.395	ng/ul	99
13) Anthracene	17.35	178	15153	0.475	ng/ul	99
15) Fluoranthene	19.28	202	20226	0.441	ng/ul	96
17) Pyrene	19.64	202	20735	0.436	ng/ul	95
18) Benzo(a)anthracene	21.39	228	19407	0.463	ng/ul	96
19) Chrysene	21.44	228	18695	0.390	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	17753	0.361	ng/ul	90
22) Benzo(k)fluoranthene	23.07	252	18269	0.381	ng/ul#	91
23) Benzo(a)pyrene	23.62	252	16978	0.399	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.09	276	19512	0.399	ng/ul	98
25) Dibenzo(a,h)anthracene	26.10	278	15912	0.405	ng/ul#	91
26) Benzo(g,h,i)perylene	26.82	276	16326	0.384	ng/ul#	91

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

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