

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004070.D
 Acq On : 18 Dec 2018 11:26
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.428

Quant Time: Dec 19 00:32:22 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 : Tue Dec 18 06:35:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7375	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4518	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	11654	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13028	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	12164	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4721	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	14733	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	8341	0.403	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	6236	0.431	ng/ul	99
7) Acenaphthylene	14.19	152	9697	0.477	ng/ul	99
8) Acenaphthene	14.53	153	7323	0.411	ng/ul	99
9) Fluorene	15.52	166	8563	0.412	ng/ul	99
11) Pentachlorophenol	16.88	266	784	0.374	ng/ul	98
12) Phenanthrene	17.26	178	15140	0.396	ng/ul	99
13) Anthracene	17.35	178	14470	0.460	ng/ul	99
15) Fluoranthene	19.28	202	19470	0.431	ng/ul	95
17) Pyrene	19.64	202	20108	0.430	ng/ul	98
18) Benzo(a)anthracene	21.39	228	19207	0.467	ng/ul	98
19) Chrysene	21.44	228	18613	0.395	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	18449	0.367	ng/ul	95
22) Benzo(k)fluoranthene	23.07	252	18601	0.379	ng/ul#	95
23) Benzo(a)pyrene	23.63	252	17496	0.402	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.09	276	20263	0.406	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.10	278	16459	0.409	ng/ul#	95
26) Benzo(g,h,i)perylene	26.82	276	16964	0.391	ng/ul	94

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

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