

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070218\
 Data File : BN001927.D
 Acq On : 02 Jul 2018 08:54
 Operator : JU/SJ
 Sample : SSTDC000.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.482

Quant Time: Jul 03 01:45:08 2018
 Quant Method : Z:\svoasrv\HPCHEM1\BNA N\Methods\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 15 13:44:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	4736	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.48	136	19858	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.34	164	11956	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.09	188	29659	0.40	ng/ul	-0.01
16) Chrysene-d12	21.28	240	24570	0.40	ng/ul	-0.01
20) Perylene-d12	23.51	264	20964	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	12991	0.41	ng/ul	-0.02
14) Fluoranthene-d10	19.12	212	33695	0.38	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.53	128	20189	0.394	ng/ul	96
5) 2-Methylnaphthalene	12.15	142	14706	0.402	ng/ul	98
7) Acenaphthylene	14.06	152	19819	0.339	ng/ul	95
8) Acenaphthene	14.40	153	16673	0.388	ng/ul	97
9) Fluorene	15.39	166	19436	0.387	ng/ul#	94
11) Pentachlorophenol	16.74	266	1791	0.319	ng/ul	99
12) Phenanthrene	17.12	178	34729	0.392	ng/ul#	88
13) Anthracene	17.22	178	29104	0.347	ng/ul#	92
15) Fluoranthene	19.15	202	39542	0.376	ng/ul	96
17) Pyrene	19.51	202	39260	0.440	ng/ul	97
18) Benzo(a)anthracene	21.27	228	30123	0.355	ng/ul#	85
19) Chrysene	21.32	228	33191	0.393	ng/ul	97
21) Benzo(b)fluoranthene	22.85	252	30135	0.396	ng/ul	86
22) Benzo(k)fluoranthene	22.89	252	31304	0.413	ng/ul#	88
23) Benzo(a)pyrene	23.41	252	26550	0.376	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.74	276	29736	0.375	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.76	278	22143	0.350	ng/ul#	91
26) Benzo(g,h,i)perylene	26.41	276	25579	0.379	ng/ul	97

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

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