

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002462.D
 Acq On : 23 Aug 2018 13:01
 Operator : JU/SJ
 Sample : SSTD1.612
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.612

Quant Time: Aug 23 14:02:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 13:07:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	7295	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4133	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	9455	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	10909	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	9687	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	18983	1.64	ng/ul	-0.02
14) Fluoranthene-d10	19.09	212	49030	1.71	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.50	128	29938	1.592	ng/ul	96
5) 2-Methylnaphthalene	12.12	142	21489	1.600	ng/ul	98
7) Acenaphthylene	14.02	152	32033	1.585	ng/ul	96
8) Acenaphthene	14.37	153	24059	1.619	ng/ul	96
9) Fluorene	15.36	166	28130	1.621	ng/ul#	94
11) Pentachlorophenol	16.72	266	2604	1.455	ng/ul	98
12) Phenanthrene	17.10	178	46973	1.665	ng/ul#	89
13) Anthracene	17.19	178	43727	1.635	ng/ul#	92
15) Fluoranthene	19.12	202	58439	1.745	ng/ul	97
17) Pyrene	19.49	202	56332	1.421	ng/ul	95
18) Benzo(a)anthracene	21.24	228	55674	1.476	ng/ul#	85
19) Chrysene	21.29	228	57680	1.537	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	57637	1.640	ng/ul	83
22) Benzo(k)fluoranthene	22.86	252	60403	1.723	ng/ul#	84
23) Benzo(a)pyrene	23.38	252	55477	1.700	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.69	276	62033	1.692	ng/ul#	85
25) Dibenzo(a,h)anthracene	25.70	278	49110	1.681	ng/ul#	87
26) Benzo(g,h,i)perylene	26.37	276	51676	1.658	ng/ul	94

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

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