

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN007022.D
 Acq On : 08 Aug 2019 12:52
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.429

Quant Time: Aug 08 13:24:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	6775	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	25664	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	15632	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	39490	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	33506	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	31428	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	17272	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	47619	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.28	128	27800	0.376	ng/ul	97
5) 2-Methylnaphthalene	11.91	142	20553	0.371	ng/ul	100
7) Acenaphthylene	13.82	152	33266	0.367	ng/ul	100
8) Acenaphthene	14.18	153	23244	0.375	ng/ul	99
9) Fluorene	15.17	166	27633	0.368	ng/ul	100
11) Pentachlorophenol	16.52	266	3674	0.338	ng/ul#	82
12) Phenanthrene	16.90	178	47084	0.374	ng/ul	99
13) Anthracene	16.99	178	46319	0.360	ng/ul	99
15) Fluoranthene	18.93	202	57636	0.368	ng/ul	98
17) Pyrene	19.29	202	58611	0.392	ng/ul	99
18) Benzo(a)anthracene	21.06	228	55002	0.374	ng/ul	99
19) Chrysene	21.11	228	51701	0.375	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	48027	0.377	ng/ul	100
22) Benzo(k)fluoranthene	22.62	252	46492	0.376	ng/ul	98
23) Benzo(a)pyrene	23.11	252	43750	0.365	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.30	276	46674	0.361	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.32	278	38220	0.370	ng/ul	98
26) Benzo(g,h,i)perylene	25.93	276	38450	0.354	ng/ul	99

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

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