

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007484.D
 Acq On : 29 Aug 2019 00:35
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.437

Quant Time: Aug 29 01:07:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 20:30:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	4224	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	18002	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	9423	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	21295	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	19898	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	20402	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	11004	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	27944	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.20	128	19825	0.390	ng/ul	99
5) 2-Methylnaphthalene	11.83	142	13471	0.395	ng/ul	98
7) Acenaphthylene	13.76	152	19638	0.375	ng/ul	99
8) Acenaphthene	14.11	153	14544	0.386	ng/ul	99
9) Fluorene	15.11	166	15916	0.390	ng/ul	97
11) Pentachlorophenol	16.46	266	2035	0.424	ng/ul	99
12) Phenanthrene	16.84	178	26470	0.398	ng/ul	100
13) Anthracene	16.93	178	25201	0.402	ng/ul	100
15) Fluoranthene	18.87	202	35396	0.398	ng/ul	100
17) Pyrene	19.24	202	36609	0.376	ng/ul	98
18) Benzo(a)anthracene	21.00	228	33822	0.380	ng/ul	99
19) Chrysene	21.05	228	36392	0.381	ng/ul	99
21) Benzo(b)fluoranthene	22.50	252	34933	0.392	ng/ul	98
22) Benzo(k)fluoranthene	22.54	252	37128	0.399	ng/ul#	91
23) Benzo(a)pyrene	23.03	252	34161	0.385	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.18	276	39158	0.394	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.19	278	30776	0.391	ng/ul	100
26) Benzo(g,h,i)perylene	25.81	276	33117	0.390	ng/ul	99

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

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