

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007459.D
 Acq On : 28 Aug 2019 07:29
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.435

Quant Time: Aug 28 08:00:28 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	3693	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	16411	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	8599	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	19016	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	15779	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	16265	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	9997	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	24301	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	17725	0.382	ng/ul	99
5) 2-Methylnaphthalene	11.84	142	12160	0.391	ng/ul	99
7) Acenaphthylene	13.76	152	18469	0.386	ng/ul	99
8) Acenaphthene	14.11	153	13267	0.386	ng/ul	99
9) Fluorene	15.11	166	14382	0.387	ng/ul	98
11) Pentachlorophenol	16.47	266	1880	0.439	ng/ul	97
12) Phenanthrene	16.84	178	23148	0.390	ng/ul	99
13) Anthracene	16.94	178	22914	0.410	ng/ul	99
15) Fluoranthene	18.87	202	30093	0.379	ng/ul	99
17) Pyrene	19.24	202	31404	0.407	ng/ul	97
18) Benzo(a)anthracene	21.01	228	27554	0.391	ng/ul	98
19) Chrysene	21.06	228	28932	0.382	ng/ul	98
21) Benzo(b)fluoranthene	22.51	252	27527	0.387	ng/ul#	77
22) Benzo(k)fluoranthene	22.55	252	28534	0.384	ng/ul#	68
23) Benzo(a)pyrene	23.04	252	27459	0.388	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.19	276	32490	0.410	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.20	278	25183	0.402	ng/ul#	83
26) Benzo(g,h,i)perylene	25.82	276	27397	0.405	ng/ul#	92

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

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