

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080319\
 Data File : BM021824.D
 Acq On : 04 Aug 2019 23:17
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.429

Quant Time: Aug 05 07:01:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 05 06:57:04 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	659	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	2894	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	1584	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3560	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3393	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4044	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1918	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	4566	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.48	128	3239	0.396	ng/ul	99
5) 2-Methylnaphthalene	12.12	142	2231	0.410	ng/ul	99
7) Acenaphthylene	14.02	152	3085	0.418	ng/ul	99
8) Acenaphthene	14.36	153	2564	0.424	ng/ul	97
9) Fluorene	15.36	166	2703	0.403	ng/ul	97
11) Pentachlorophenol	16.74	266	282	0.429	ng/ul	86
12) Phenanthrene	17.10	178	4174	0.403	ng/ul	99
13) Anthracene	17.20	178	3496	0.396	ng/ul	99
15) Fluoranthene	19.12	202	5253	0.379	ng/ul	98
17) Pyrene	19.48	202	5533	0.397	ng/ul#	94
18) Benzo(a)anthracene	21.24	228	4799	0.395	ng/ul	100
19) Chrysene	21.28	228	6299	0.404	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	5570	0.408	ng/ul	96
22) Benzo(k)fluoranthene	22.85	252	6154	0.399	ng/ul	98
23) Benzo(a)pyrene	23.37	252	5564	0.401	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.70	276	6728	0.406	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.71	278	5386	0.402	ng/ul	96
26) Benzo(g,h,i)perylene	26.37	276	5998	0.402	ng/ul	98

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

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