

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091719\
 Data File : BM022709.D
 Acq On : 17 Sep 2019 17:28
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.469

Quant Time: Sep 17 18:35:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	8639	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	34762	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	18740	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	45067	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	37604	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	32313	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	21531	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	47654	0.39	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.68	128	36181	0.365	ng/ul 100
5) 2-Methylnaphthalene	12.30	142	24943	0.367	ng/ul 100
7) Acenaphthylene	14.19	152	31343	0.361	ng/ul 99
8) Acenaphthene	14.53	153	29081	0.400	ng/ul 100
9) Fluorene	15.52	166	30491	0.366	ng/ul 100
11) Pentachlorophenol	16.86	266	3260	0.357	ng/ul 99
12) Phenanthrene	17.25	178	51744	0.360	ng/ul 99
13) Anthracene	17.34	178	43067	0.350	ng/ul 100
15) Fluoranthene	19.27	202	57961	0.374	ng/ul 99
17) Pyrene	19.63	202	62612	0.367	ng/ul 99
18) Benzo(a)anthracene	21.37	228	49617	0.345	ng/ul 99
19) Chrysene	21.43	228	54890	0.363	ng/ul 100
21) Benzo(b)fluoranthene	22.98	252	52219	0.372	ng/ul 99
22) Benzo(k)fluoranthene	23.03	252	51518	0.360	ng/ul 98
23) Benzo(a)pyrene	23.57	252	44466	0.362	ng/ul 98
24) Indeno(1,2,3-cd)pyrene	25.98	276	49813	0.332	ng/ul# 99
25) Dibenzo(a,h)anthracene	26.00	278	41039	0.339	ng/ul 99
26) Benzo(g,h,i)perylene	26.69	276	41879	0.351	ng/ul 99

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

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