

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030317\
 Data File : BM009072.D
 Acq On : 03 Mar 2017 17:35
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV40

Quant Time: Mar 03 23:59:24 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 03 23:57:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	142	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	397	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	218	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.28	188	513	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	568	0.40	ng/ul	0.00
20) Perylene-d12	23.79	264	525	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	255	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	635	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.74	128	440	0.38	ng/ul#	74
5) 2-Methylnaphthalene	12.35	142	325	0.39	ng/ul	94
7) Acenaphthylene	14.25	152	461	0.41	ng/ul#	80
8) Acenaphthene	14.59	153	360	0.41	ng/ul	96
9) Fluorene	15.58	166	413	0.41	ng/ul#	93
11) Pentachlorophenol	16.92	266	99	0.40	ng/ul	89
12) Phenanthrene	17.31	178	634	0.39	ng/ul#	89
13) Anthracene	17.40	178	632	0.40	ng/ul#	87
15) Fluoranthene	19.33	202	791	0.38	ng/ul	93
17) Pyrene	19.69	202	820	0.42	ng/ul#	91
18) Benzo(a)anthracene	21.43	228	763	0.40	ng/ul#	93
19) Chrysene	21.48	228	729	0.41	ng/ul#	91
21) Benzo(b)fluoranthene	23.07	252	770	0.37	ng/ul	89
22) Benzo(k)fluoranthene	23.12	252	807	0.44	ng/ul#	85
23) Benzo(a)pyrene	23.67	252	767	0.41	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.16	276	883	0.37	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	741	0.37	ng/ul#	88
26) Benzo(g,h,i)perylene	26.91	276	758	0.38	ng/ul#	88

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

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