

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060717\
 Data File : BE093071.D
 Acq On : 7 Jun 2017 16:30
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV58

Quant Time: Jun 07 17:12:37 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 07 16:36:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	4540	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	20850	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10743	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	30199	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	37044	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	30526	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	12153	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	33199	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.57	128	20822	0.40	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	14005	0.40	ng/ul	99
7) Acenaphthylene	14.09	152	18289	0.40	ng/ul	95
8) Acenaphthene	14.43	153	15192	0.41	ng/ul	98
9) Fluorene	15.42	166	17713	0.41	ng/ul#	91
11) Pentachlorophenol	16.75	266	1529	0.34	ng/ul	97
12) Phenanthrene	17.13	178	33579	0.40	ng/ul#	86
13) Anthracene	17.22	178	28856	0.40	ng/ul#	86
15) Fluoranthene	19.16	202	41045	0.40	ng/ul	83
17) Pyrene	19.51	202	45514	0.42	ng/ul#	87
18) Benzo(a)anthracene	21.25	228	39792	0.40	ng/ul#	85
19) Chrysene	21.30	228	41586	0.41	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	36431	0.39	ng/ul	86
22) Benzo(k)fluoranthene	22.98	252	39559	0.40	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	34792	0.40	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.19	276	39664	0.39	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	33352	0.39	ng/ul#	91
26) Benzo(g,h,i)perylene	26.96	276	34251	0.39	ng/ul	94

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

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