

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE121817\
 Data File : BE095010.D
 Acq On : 18 Dec 2017 15:13
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV07

Quant Time: Dec 18 16:18:01 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:40:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4667	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14590	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	7977	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	19468	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	19824	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	16332	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	9616	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	24047	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.34	128	15212	0.393	ng/ul#	88
5) 2-Methylnaphthalene	12.91	142	10675	0.355	ng/ul	100
7) Acenaphthylene	14.77	152	14440	0.396	ng/ul	94
8) Acenaphthene	15.11	153	11119	0.397	ng/ul	95
9) Fluorene	16.07	166	11953	0.384	ng/ul	91
11) Pentachlorophenol	17.41	266	647	0.327	ng/ul#	59
12) Phenanthrene	17.78	178	19844	0.387	ng/ul#	81
13) Anthracene	17.87	178	17596	0.362	ng/ul#	86
15) Fluoranthene	19.76	202	26274	0.393	ng/ul	85
17) Pyrene	20.11	202	26360	0.383	ng/ul#	82
18) Benzo(a)anthracene	21.82	228	22230	0.390	ng/ul#	85
19) Chrysene	21.88	228	23986	0.397	ng/ul	94
21) Benzo(b)fluoranthene	23.67	252	22802	0.400	ng/ul	85
22) Benzo(k)fluoranthene	23.73	252	20275	0.390	ng/ul#	85
23) Benzo(a)pyrene	24.38	252	19690	0.402	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.34	276	21305	0.396	ng/ul#	82
25) Dibenzo(a,h)anthracene	27.36	278	17338	0.398	ng/ul#	89
26) Benzo(g,h,i)perylene	28.22	276	18351	0.394	ng/ul	95

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

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