

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097296.D
 Acq On : 17 Aug 2018 14:09
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV07

Quant Time: Aug 17 14:55:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 14:15:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	944	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4782	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2440	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	5570	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	4784	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	3998	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	3016	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5798	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	4670	0.400	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	3002	0.399	ng/ul	99
7) Acenaphthylene	13.73	152	4383	0.398	ng/ul	98
8) Acenaphthene	14.08	153	3161	0.399	ng/ul#	88
9) Fluorene	15.08	166	3423	0.397	ng/ul	91
11) Pentachlorophenol	16.44	266	115	0.318	ng/ul	95
12) Phenanthrene	16.81	178	5732	0.398	ng/ul#	89
13) Anthracene	16.90	178	4737	0.393	ng/ul#	92
15) Fluoranthene	18.84	202	6052	0.396	ng/ul	85
17) Pyrene	19.20	202	5871	0.406	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	4611	0.389	ng/ul#	86
19) Chrysene	21.01	228	5705	0.403	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	4277	0.395	ng/ul	85
22) Benzo(k)fluoranthene	22.58	252	5461	0.406	ng/ul#	88
23) Benzo(a)pyrene	23.12	252	3841	0.384	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.54	276	4925	0.396	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.55	278	3994	0.397	ng/ul	95
26) Benzo(g,h,i)perylene	26.22	276	3882	0.390	ng/ul	93

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

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