

Data Path : U:\HPCHEM1\BNA E\DATA\BE012618\
 Data File : BE095125.D
 Acq On : 26 Jan 2018 13:36
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV09

Quant Time: Jan 26 14:12:31 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 26 13:41:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2927	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8315	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4425	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	10540	0.40	ng/ul	0.02
16) Chrysene-d12	21.81	240	14374	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	12624	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	5451	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	15765	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	9225	0.396	ng/ul#	90
5) 2-Methylnaphthalene	12.89	142	6235	0.401	ng/ul	99
7) Acenaphthylene	14.74	152	8475	0.398	ng/ul	96
8) Acenaphthene	15.08	153	6607	0.400	ng/ul	97
9) Fluorene	16.02	166	6754	0.397	ng/ul#	70
11) Pentachlorophenol	17.36	266	644	0.315	ng/ul	95
12) Phenanthrene	17.75	178	13815	0.390	ng/ul#	85
13) Anthracene	17.84	178	12725	0.392	ng/ul#	90
15) Fluoranthene	19.73	202	18063	0.394	ng/ul	83
17) Pyrene	20.08	202	19883	0.412	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	17537	0.390	ng/ul#	85
19) Chrysene	21.84	228	18792	0.403	ng/ul	95
21) Benzo(b)fluoranthene	23.63	252	17108	0.384	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	17281	0.424	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	15634	0.393	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.25	276	17736	0.407	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	14381	0.403	ng/ul#	87
26) Benzo(g,h,i)perylene	28.12	276	14988	0.406	ng/ul#	94

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

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