

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040618\
 Data File : BE095959.D
 Acq On : 6 Apr 2018 16:39
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV35

Quant Time: Apr 06 23:43:18 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 06 17:33:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1818	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4610	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.98	164	2605	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6390	0.40	ng/ul	0.00
16) Chrysene-d12	21.78	240	8483	0.40	ng/ul	0.00
20) Perylene-d12	24.40	264	7046	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	3374	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.67	212	10214	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	4992	0.388	ng/ul#	89
5) 2-Methylnaphthalene	12.84	142	3397	0.389	ng/ul	99
7) Acenaphthylene	14.70	152	5195	0.385	ng/ul	97
8) Acenaphthene	15.04	153	3767	0.389	ng/ul	94
9) Fluorene	16.00	166	4269	0.390	ng/ul#	90
11) Pentachlorophenol	17.34	266	450	0.333	ng/ul	97
12) Phenanthrene	17.73	178	7296	0.387	ng/ul#	89
13) Anthracene	17.82	178	7190	0.384	ng/ul#	91
15) Fluoranthene	19.69	202	9796	0.382	ng/ul	84
17) Pyrene	20.05	202	10776	0.374	ng/ul#	78
18) Benzo(a)anthracene	21.76	228	12841	0.375	ng/ul#	85
19) Chrysene	21.82	228	9277	0.376	ng/ul	96
21) Benzo(b)fluoranthene	23.59	252	8781	0.385	ng/ul	86
22) Benzo(k)fluoranthene	23.64	252	8803	0.388	ng/ul#	89
23) Benzo(a)pyrene	24.28	252	8515	0.387	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.20	276	9588	0.392	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.22	278	7858	0.399	ng/ul	90
26) Benzo(g,h,i)perylene	28.07	276	7964	0.390	ng/ul	96

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

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