

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096959.D
 Acq On : 17 Jul 2018 14:51
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV67

Quant Time: Jul 17 15:29:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 14:56:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1580	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	7083	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3513	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	9834	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12757	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11713	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4423	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	11833	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.19	128	6924	0.403	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	4496	0.397	ng/ul	98
7) Acenaphthylene	13.74	152	6143	0.401	ng/ul	98
8) Acenaphthene	14.08	153	4755	0.400	ng/ul#	93
9) Fluorene	15.08	166	4945	0.392	ng/ul	94
11) Pentachlorophenol	16.43	266	368	0.331	ng/ul	95
12) Phenanthrene	16.82	178	10180	0.394	ng/ul#	89
13) Anthracene	16.90	178	8775	0.389	ng/ul#	90
15) Fluoranthene	18.84	202	13296	0.391	ng/ul	83
17) Pyrene	19.21	202	14238	0.386	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	12992	0.385	ng/ul#	86
19) Chrysene	21.01	228	15034	0.392	ng/ul	95
21) Benzo(b)fluoranthene	22.54	252	12024	0.362	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	13914	0.405	ng/ul#	84
23) Benzo(a)pyrene	23.11	252	10772	0.369	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.50	276	12458	0.381	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	10474	0.383	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	11254	0.386	ng/ul	94

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

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