

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097209.D
 Acq On : 4 Aug 2018 19:52
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.489

Quant Time: Aug 06 03:29:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 06 00:53:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1437	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7946	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4709	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11005	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11489	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11851	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	5171	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12890	0.34	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.18	128	7422	0.386	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	5126	0.403	ng/ul	98
7) Acenaphthylene	13.73	152	7997	0.393	ng/ul	97
8) Acenaphthene	14.08	153	5705	0.387	ng/ul	93
9) Fluorene	15.08	166	6495	0.384	ng/ul#	91
11) Pentachlorophenol	16.44	266	543	0.344	ng/ul	97
12) Phenanthrene	16.81	178	10831	0.386	ng/ul#	88
13) Anthracene	16.90	178	10526	0.402	ng/ul#	92
15) Fluoranthene	18.84	202	12715	0.343	ng/ul	84
17) Pyrene	19.20	202	12770	0.415	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	13210	0.402	ng/ul#	84
19) Chrysene	21.01	228	13239	0.391	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	12709	0.393	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	13744	0.395	ng/ul#	89
23) Benzo(a)pyrene	23.11	252	12928	0.400	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.51	276	16180	0.400	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.53	278	13462	0.407	ng/ul#	90
26) Benzo(g,h,i)perylene	26.21	276	13524	0.387	ng/ul#	92

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

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