

Data Path : \\Terastorage\svoasrv\HPCHEM1\BNA_E\Data\BE012618\
 Data File : BE095124.D
 Acq On : 26 Jan 2018 13:00
 Operator : SJ/JU
 Sample : SSTD3.208
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD3.208

Quant Time: Jan 26 13:40:33 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 12:06:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2938	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8573	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4800	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	10688	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	15424	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	12963	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	46845	3.32	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	142777	4.30	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.32	128	76730	3.378	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	53628	3.388	ng/ul	100
7) Acenaphthylene	14.74	152	79838	3.713	ng/ul	95
8) Acenaphthene	15.08	153	58642	3.506	ng/ul	96
9) Fluorene	16.02	166	62107	3.480	ng/ul#	69
11) Pentachlorophenol	17.36	266	8039	12.717	ng/ul	97
12) Phenanthrene	17.75	178	119823	4.222	ng/ul#	85
13) Anthracene	17.85	178	121878	4.923	ng/ul#	90
15) Fluoranthene	19.73	202	162823	4.489	ng/ul	82
17) Pyrene	20.08	202	166929	2.977	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	162435	3.806	ng/ul#	85
19) Chrysene	21.84	228	161242	3.386	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	148433	3.401	ng/ul	81
22) Benzo(k)fluoranthene	23.67	252	154836	3.846	ng/ul#	84
23) Benzo(a)pyrene	24.32	252	141487	3.797	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	27.25	276	156495	4.730	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	128897	5.454	ng/ul#	85
26) Benzo(g,h,i)perylene	28.12	276	129437	3.934	ng/ul#	91

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

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