

Data Path : Z:\HPCHEM1\BNA_E\Data\BE121817\
 Data File : BE095009.D
 Acq On : 18 Dec 2017 14:37
 Operator : SJ/JU
 Sample : SSTD3.206
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD3.206

Quant Time: Dec 18 15:37:09 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:33:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	4775	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	15681	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8923	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	20880	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	22991	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	18054	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	85596	3.38	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	230614	3.21	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	132505	3.548	ng/ul#	88
5) 2-Methylnaphthalene	12.91	142	96996	3.734	ng/ul	100
7) Acenaphthylene	14.77	152	143992	3.524	ng/ul	95
8) Acenaphthene	15.11	153	102740	3.572	ng/ul	94
9) Fluorene	16.07	166	116274	3.333	ng/ul	92
11) Pentachlorophenol	17.41	266	9150	8.311	ng/ul#	60
12) Phenanthrene	17.78	178	175471	3.129	ng/ul#	80
13) Anthracene	17.87	178	176088	3.082	ng/ul#	85
15) Fluoranthene	19.76	202	252774	3.435	ng/ul	84
17) Pyrene	20.11	202	253687	3.526	ng/ul#	81
18) Benzo(a)anthracene	21.82	228	225167	3.221	ng/ul#	84
19) Chrysene	21.89	228	220504	3.160	ng/ul	98
21) Benzo(b)fluoranthene	23.68	252	212184	3.945	ng/ul	82
22) Benzo(k)fluoranthene	23.73	252	196934	3.443	ng/ul#	83
23) Benzo(a)pyrene	24.38	252	189248	3.538	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.34	276	209411	3.592	ng/ul#	82
25) Dibenzo(a,h)anthracene	27.36	278	174193	3.545	ng/ul#	86
26) Benzo(g,h,i)perylene	28.22	276	174563	3.573	ng/ul	94

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

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