

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102017\
 Data File : BE094482.D
 Acq On : 20 Oct 2017 19:42
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.445

Quant Time: Oct 22 02:48:06 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 18 17:51:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1362	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7727	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	4513	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10327	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	10849	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	10600	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4938	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	12443	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	7480	0.395	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	5265	0.429	ng/ul	98
7) Acenaphthylene	14.24	152	8407	0.478	ng/ul	97
8) Acenaphthene	14.58	153	5880	0.406	ng/ul	98
9) Fluorene	15.58	166	6705	0.395	ng/ul	92
11) Pentachlorophenol	16.98	266	426	0.352	ng/ul	94
12) Phenanthrene	17.29	178	11112	0.412	ng/ul#	86
13) Anthracene	17.39	178	10802	0.412	ng/ul#	92
15) Fluoranthene	19.31	202	12793	0.361	ng/ul	85
17) Pyrene	19.67	202	13596	0.450	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	12768	0.444	ng/ul#	86
19) Chrysene	21.46	228	12392	0.373	ng/ul	98
21) Benzo(b)fluoranthene	23.17	252	11999	0.417	ng/ul	89
22) Benzo(k)fluoranthene	23.22	252	12283	0.369	ng/ul	91
23) Benzo(a)pyrene	23.83	252	11547	0.386	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.63	276	12750	0.417	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.63	278	10860	0.423	ng/ul	94
26) Benzo(g,h,i)perylene	27.47	276	10035	0.380	ng/ul	97

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

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