

Data Path : Z:\HPCHEM1\BNA_E\Data\BE063017\
 Data File : BE093373.D
 Acq On : 30 Jun 2017 8:59
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.427

Quant Time: Jun 30 10:41:20 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 30 10:40:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	4380	0.40	ng/ul	0.00
2) Naphthalene-d8	10.74	136	21301	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	12935	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	36197	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	50856	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	48862	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.32	152	13287	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	46058	0.47	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.79	128	21703	0.411	ng/ul#	91
5) 2-Methylnaphthalene	12.39	142	15665	0.433	ng/ul	97
7) Acenaphthylene	14.28	152	23342	0.421	ng/ul	95
8) Acenaphthene	14.62	153	17675	0.395	ng/ul	97
9) Fluorene	15.60	166	21890	0.417	ng/ul#	92
11) Pentachlorophenol	16.94	266	1257	0.234	ng/ul	95
12) Phenanthrene	17.33	178	40321	0.405	ng/ul#	92
13) Anthracene	17.41	178	39285	0.453	ng/ul#	91
15) Fluoranthene	19.33	202	52376	0.429	ng/ul	85
17) Pyrene	19.69	202	56018	0.374	ng/ul#	84
18) Benzo(a)anthracene	21.42	228	57752	0.426	ng/ul#	85
19) Chrysene	21.48	228	56469	0.401	ng/ul	97
21) Benzo(b)fluoranthene	23.19	252	57839	0.390	ng/ul	84
22) Benzo(k)fluoranthene	23.24	252	62060	0.394	ng/ul#	89
23) Benzo(a)pyrene	23.85	252	55786	0.398	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.62	276	64649	0.398	ng/ul#	82
25) Dibenzo(a,h)anthracene	26.65	278	54935	0.402	ng/ul#	87
26) Benzo(g,h,i)perylene	27.43	276	55315	0.397	ng/ul	94

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

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