

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020118\
 Data File : BE095373.D
 Acq On : 2 Feb 2018 00:41
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.432

Quant Time: Feb 02 04:09:46 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 Feb 02 04:07:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2418	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7417	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4276	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8761	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9576	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9096	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5113	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	12417	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8312	0.400	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5915	0.427	ng/ul	100
7) Acenaphthylene	14.74	152	8860	0.431	ng/ul	97
8) Acenaphthene	15.08	153	6423	0.402	ng/ul	95
9) Fluorene	16.02	166	6869	0.417	ng/ul#	79
11) Pentachlorophenol	17.36	266	635	0.374	ng/ul	96
12) Phenanthrene	17.75	178	12067	0.410	ng/ul#	89
13) Anthracene	17.85	178	11567	0.429	ng/ul#	92
15) Fluoranthene	19.72	202	13860	0.363	ng/ul	83
17) Pyrene	20.07	202	15291	0.476	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	12948	0.432	ng/ul#	86
19) Chrysene	21.84	228	12425	0.400	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	12164	0.379	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	11923	0.406	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	11612	0.405	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.25	276	12728	0.405	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.27	278	10232	0.398	ng/ul#	89
26) Benzo(g,h,i)perylene	28.12	276	10712	0.403	ng/ul#	93

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

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