

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095431.D
 Acq On : 3 Feb 2018 11:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.437

Quant Time: Feb 05 01:50:59 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 01:47:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	2642	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	8182	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4660	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10540	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	11081	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	10398	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5544	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	14155	0.35	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.31	128	9177	0.401	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	6327	0.414	ng/ul	99
7) Acenaphthylene	14.74	152	9601	0.428	ng/ul	97
8) Acenaphthene	15.08	153	7085	0.407	ng/ul	96
9) Fluorene	16.02	166	7964	0.444	ng/ul#	84
11) Pentachlorophenol	17.36	266	814	0.399	ng/ul	97
12) Phenanthrene	17.75	178	13682	0.386	ng/ul#	90
13) Anthracene	17.84	178	12613	0.389	ng/ul#	92
15) Fluoranthene	19.72	202	15916	0.347	ng/ul	82
17) Pyrene	20.07	202	17422	0.469	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	14852	0.429	ng/ul#	84
19) Chrysene	21.84	228	14406	0.400	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	14410	0.393	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	13270	0.395	ng/ul#	85
23) Benzo(a)pyrene	24.31	252	13358	0.408	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.24	276	14465	0.403	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.26	278	11486	0.391	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	12158	0.400	ng/ul#	93

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

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