

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095870.D
 Acq On : 28 Mar 2018 10:38
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.423

Quant Time: Mar 28 16:21:58 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 16 02:40:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1580	0.40	ng/ul	-0.03
2) Naphthalene-d8	11.24	136	4100	0.40	ng/ul	-0.04
6) Acenaphthene-d10	15.00	164	2546	0.40	ng/ul	-0.04
10) Phenanthrene-d10	17.70	188	6647	0.40	ng/ul	-0.04
16) Chrysene-d12	21.79	240	7608	0.40	ng/ul	-0.05
20) Perylene-d12	24.42	264	7241	0.40	ng/ul	-0.08
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2951	0.42	ng/ul	-0.03
14) Fluoranthene-d10	19.68	212	10150	0.48	ng/ul	-0.04
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	4246	0.371	ng/ul#	90
5) 2-Methylnaphthalene	12.87	142	2958	0.372	ng/ul	99
7) Acenaphthylene	14.72	152	4740	0.369	ng/ul	98
8) Acenaphthene	15.06	153	3434	0.355	ng/ul	96
9) Fluorene	16.02	166	4007	0.381	ng/ul#	89
11) Pentachlorophenol	17.36	266	348	0.372	ng/ul	87
12) Phenanthrene	17.75	178	7066	0.355	ng/ul#	90
13) Anthracene	17.83	178	6907	0.371	ng/ul#	90
15) Fluoranthene	19.71	202	9620	0.399	ng/ul	84
17) Pyrene	20.06	202	10852	0.376	ng/ul#	78
18) Benzo(a)anthracene	21.77	228	9383	0.382	ng/ul#	86
19) Chrysene	21.83	228	8850	0.354	ng/ul	97
21) Benzo(b)fluoranthene	23.60	252	8582	0.342	ng/ul	86
22) Benzo(k)fluoranthene	23.65	252	8621	0.362	ng/ul#	88
23) Benzo(a)pyrene	24.30	252	8237	0.360	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.22	276	9336	0.368	ng/ul#	76
25) Dibenzo(a,h)anthracene	27.24	278	7553	0.372	ng/ul	92
26) Benzo(g,h,i)perylene	28.09	276	7771	0.356	ng/ul	96

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

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