

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096173.D
 Acq On : 26 Apr 2018 10:21
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.465

Quant Time: Apr 27 02:25:56 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 26 06:40:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	667	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2477	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1456	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3616	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	5026	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3858	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1856	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5969	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	2876	0.417	ng/ul#	91
5) 2-Methylnaphthalene	12.81	142	1876	0.406	ng/ul	98
7) Acenaphthylene	14.67	152	2927	0.390	ng/ul	98
8) Acenaphthene	15.01	153	2232	0.412	ng/ul	94
9) Fluorene	15.97	166	2950	0.412	ng/ul	90
11) Pentachlorophenol	17.32	266	262	0.301	ng/ul	96
12) Phenanthrene	17.70	178	5722	0.537	ng/ul#	88
13) Anthracene	17.79	178	4470	0.419	ng/ul#	91
15) Fluoranthene	19.67	202	8026	0.493	ng/ul	84
17) Pyrene	20.02	202	4778	0.474	ng/ul#	80
18) Benzo(a)anthracene	21.73	228	6179	0.406	ng/ul#	85
19) Chrysene	21.79	228	6237	0.446	ng/ul	97
21) Benzo(b)fluoranthene	23.55	252	5883	0.403	ng/ul	88
22) Benzo(k)fluoranthene	23.60	252	5389	0.386	ng/ul#	89
23) Benzo(a)pyrene	24.24	252	5165	0.407	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.13	276	5581	0.326	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.15	278	4518	0.316	ng/ul	92
26) Benzo(g,h,i)perylene	28.01	276	4635	0.311	ng/ul	95

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

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