

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040618\
 Data File : BE095957.D
 Acq On : 6 Apr 2018 15:27
 Operator : SJ/JU
 Sample : SSTD1.633
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.633

Quant Time: Apr 06 16:44:58 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 06 15:00:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1792	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4597	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.98	164	2579	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6057	0.40	ng/ul	0.00
16) Chrysene-d12	21.78	240	7190	0.40	ng/ul	0.00
20) Perylene-d12	24.40	264	7077	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	13446	1.73	ng/ul	0.00
14) Fluoranthene-d10	19.67	212	39597	2.06	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	19866	1.548	ng/ul#	88
5) 2-Methylnaphthalene	12.84	142	13596	1.524	ng/ul	99
7) Acenaphthylene	14.70	152	21256	1.632	ng/ul	96
8) Acenaphthene	15.04	153	15017	1.531	ng/ul	94
9) Fluorene	16.01	166	16893	1.587	ng/ul#	92
11) Pentachlorophenol	17.34	266	2159	2.536	ng/ul	99
12) Phenanthrene	17.72	178	28031	1.545	ng/ul#	88
13) Anthracene	17.82	178	28587	1.685	ng/ul#	91
15) Fluoranthene	19.69	202	38223	1.740	ng/ul	84
17) Pyrene	20.05	202	40035	1.466	ng/ul#	77
18) Benzo(a)anthracene	21.76	228	50093	2.159	ng/ul#	85
19) Chrysene	21.81	228	35956	1.522	ng/ul	96
21) Benzo(b)fluoranthene	23.59	252	34858	1.422	ng/ul	83
22) Benzo(k)fluoranthene	23.64	252	35538	1.525	ng/ul#	84
23) Benzo(a)pyrene	24.28	252	33975	1.518	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.20	276	38506	1.552	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.21	278	31515	1.588	ng/ul#	86
26) Benzo(g,h,i)perylene	28.07	276	32083	1.505	ng/ul#	92

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

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