

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082516\
 Data File : BE092276.D
 Acq On : 25 Aug 2016 17:30
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Aug 26 04:20:12 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE082516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 04:20:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	6099	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.49	136	27365	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	15092	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	38288	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	49194	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	46377	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	15076	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	41296	0.47	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	10.53	128	25508	0.38	ng/ul	97
5) 2-Methylnaphthalene	12.15	142	17447	0.38	ng/ul	99
7) Acenaphthylene	14.05	152	27505	0.35	ng/ul	99
8) Acenaphthene	14.39	153	18957	0.37	ng/ul	98
9) Fluorene	15.38	166	22522	0.39	ng/ul	99
11) Pentachlorophenol	16.74	266	2049	0.24	ng/ul	96
12) Phenanthrene	17.10	178	40826	0.39	ng/ul	99
13) Anthracene	17.20	178	37747	0.38	ng/ul	99
15) Fluoranthene	19.12	202	48464	0.39	ng/ul	98
17) Pyrene	19.49	202	50222	0.33	ng/ul	100
18) Benzo(a)anthracene	21.22	228	47995	0.33	ng/ul	99
19) Chrysene	21.27	228	55312	0.38	ng/ul	99
21) Benzo(b)fluoranthene	22.89	252	52798	0.33	ng/ul	99
22) Benzo(k)fluoranthene	22.94	252	57268	0.41	ng/ul	99
23) Benzo(a)pyrene	23.52	252	53854	0.38	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.14	276	60074	0.35	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.15	278	48302	0.35	ng/ul	99
26) Benzo(g,h,i)perylene	26.90	276	52206	0.35	ng/ul	99

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

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