

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082516\
 Data File : BE092279.D
 Acq On : 25 Aug 2016 19:17
 Operator : UM/SJ
 Sample : SSTD0.469
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.469

Quant Time: Aug 26 04:20:40 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	7.72	152	4960	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	22797	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	12360	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	32249	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	39104	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	36153	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	12288	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	33329	0.45	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.54	128	21583	0.38	ng/ul	100
5) 2-Methylnaphthalene	12.16	142	14267	0.38	ng/ul	100
7) Acenaphthylene	14.05	152	22323	0.35	ng/ul	100
8) Acenaphthene	14.39	153	15680	0.37	ng/ul	100
9) Fluorene	15.38	166	18352	0.38	ng/ul	100
11) Pentachlorophenol	16.74	266	1376	0.19	ng/ul	99
12) Phenanthrene	17.10	178	33701	0.38	ng/ul	100
13) Anthracene	17.20	178	29716	0.35	ng/ul	100
15) Fluoranthene	19.13	202	39078	0.38	ng/ul	100
17) Pyrene	19.49	202	41222	0.34	ng/ul	100
18) Benzo(a)anthracene	21.22	228	37394	0.32	ng/ul	100
19) Chrysene	21.27	228	44242	0.39	ng/ul	100
21) Benzo(b)fluoranthene	22.89	252	39210	0.31	ng/ul	100
22) Benzo(k)fluoranthene	22.94	252	47568	0.43	ng/ul	100
23) Benzo(a)pyrene	23.52	252	42490	0.38	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.13	276	45769	0.34	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.15	278	36320	0.34	ng/ul	100
26) Benzo(g,h,i)perylene	26.90	276	40546	0.35	ng/ul	100

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

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