

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122016\
 Data File : BM008490.D
 Acq On : 20 Dec 2016 21:42
 Operator : UM/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDC0.451

Manual Integrations
 APPROVED

sohil
 12/21/2016 6:21:49 PM

Quant Time: Dec 21 12:51:37 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 02:43:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	311	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	1132	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	625	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	1133	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	1046	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	1051	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	648	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	1199m	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.52	128	1256	0.39	ng/ul#	80
5) 2-Methylnaphthalene	12.15	142	798	0.37	ng/ul	95
7) Acenaphthylene	14.02	152	1181	0.42	ng/ul#	86
8) Acenaphthene	14.37	153	858	0.40	ng/ul	93
9) Fluorene	15.39	166	941	0.39	ng/ul#	93
11) Pentachlorophenol	16.76	266	157	0.45	ng/ul	96
12) Phenanthrene	17.11	178	1331	0.38	ng/ul#	91
13) Anthracene	17.21	178	1330	0.40	ng/ul#	89
15) Fluoranthene	19.13	202	1738	0.41	ng/ul	92
17) Pyrene	19.50	202	1764	0.44	ng/ul#	95
18) Benzo(a)anthracene	21.24	228	1387	0.40	ng/ul#	92
19) Chrysene	21.30	228	1506m	0.40	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	1436	0.35	ng/ul	94
22) Benzo(k)fluoranthene	22.86	252	1678	0.42	ng/ul#	91
23) Benzo(a)pyrene	23.39	252	1544	0.40	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.73	276	1893	0.41	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.74	278	1503	0.41	ng/ul#	87
26) Benzo(g,h,i)perylene	26.42	276	1676	0.43	ng/ul#	94

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

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