

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097290.D
 Acq On : 17 Aug 2018 10:22
 Operator : SJ/JU
 Sample : SSTDO.101
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDO.101

Manual Integrations
 APPROVED

Sohil
 8/20/2018 4:15:28 PM

Quant Time: Aug 17 12:23:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 12:19:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	918	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4706	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2448	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	5601	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	4547	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4320	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	741	0.10	ng/ul	0.01
14) Fluoranthene-d10	18.81	212	1378	0.07	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	1146	0.101	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	684	0.091	ng/ul	99
7) Acenaphthylene	13.74	152	1094	0.103	ng/ul	93
8) Acenaphthene	14.09	153	777	0.102	ng/ul#	89
9) Fluorene	15.08	166	822	0.093	ng/ul	91
12) Phenanthrene	16.81	178	1402	0.098	ng/ul#	90
13) Anthracene	16.90	178	1075	0.081	ng/ul#	93
15) Fluoranthene	18.84	202	1373	0.073	ng/ul	86
17) Pyrene	19.21	202	1500	0.123	ng/ul#	86
18) Benzo(a)anthracene	20.96	228	1081	0.083	ng/ul#	87
19) Chrysene	21.01	228	1351m	0.101	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	979	0.083	ng/ul	96
22) Benzo(k)fluoranthene	22.58	252	1491m	0.117	ng/ul	
23) Benzo(a)pyrene	23.12	252	1005	0.085	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.54	276	1251	0.085	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.55	278	974	0.081	ng/ul#	76
26) Benzo(a,h,i)perylene	26.22	276	1009	0.079	ng/ul#	82

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

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