

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097329.D
 Acq On : 18 Aug 2018 12:01
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.410

Quant Time: Aug 20 01:00:33 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 : Sat Aug 18 05:45:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	843	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	3520	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	1097	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	2820	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	2619	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	2329	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1497	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	2426	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	3132	0.365	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	1504	0.272	ng/ul	100
7) Acenaphthylene	13.73	152	1858	0.375	ng/ul	99
8) Acenaphthene	14.08	153	1553	0.436	ng/ul#	69
9) Fluorene	15.08	166	1559	0.402	ng/ul	89
11) Pentachlorophenol	16.45	266	50	0.273	ng/ul#	78
12) Phenanthrene	16.81	178	2805	0.385	ng/ul#	88
13) Anthracene	16.90	178	1953	0.320	ng/ul	92
15) Fluoranthene	18.84	202	2601	0.336	ng/ul	86
17) Pyrene	19.21	202	2084	0.263	ng/ul#	85
18) Benzo(a)anthracene	20.96	228	2165	0.333	ng/ul#	86
19) Chrysene	21.01	228	2874	0.371	ng/ul	95
21) Benzo(b)fluoranthene	22.53	252	2170	0.344	ng/ul	90
22) Benzo(k)fluoranthene	22.58	252	3212	0.410	ng/ul	92
23) Benzo(a)pyrene	23.11	252	2024	0.347	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.53	276	2967	0.409	ng/ul#	74
25) Dibenzo(a,h)anthracene	25.54	278	2458	0.419	ng/ul	97
26) Benzo(g,h,i)perylene	26.22	276	2158	0.372	ng/ul#	95

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

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