

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098135.D
 Acq On : 1 Nov 2018 2:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.478

Quant Time: Nov 01 04:21:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 02:11:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1213	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	6196	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4234	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	9957	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	12885	0.40	ng/ul	0.00
20) Perylene-d12	24.01	264	11131	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4203	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	12578	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.72	128	5804	0.391	ng/ul	98
5) 2-Methylnaphthalene	12.34	142	4186	0.373	ng/ul	97
7) Acenaphthylene	14.24	152	6460	0.387	ng/ul	98
8) Acenaphthene	14.59	153	4709	0.377	ng/ul	99
9) Fluorene	15.58	166	5697	0.379	ng/ul	98
11) Pentachlorophenol	16.94	266	530	1.017	ng/ul	92
12) Phenanthrene	17.32	178	8991	0.391	ng/ul	99
13) Anthracene	17.42	178	8997	0.396	ng/ul	99
15) Fluoranthene	19.35	202	11381	0.388	ng/ul	98
17) Pyrene	19.71	202	13108	0.436	ng/ul	97
18) Benzo(a)anthracene	21.45	228	17882	0.472	ng/ul	99
19) Chrysene	21.50	228	12099	0.413	ng/ul	100
21) Benzo(b)fluoranthene	23.23	252	11950	0.387	ng/ul	97
22) Benzo(k)fluoranthene	23.28	252	12175	0.403	ng/ul#	98
23) Benzo(a)pyrene	23.90	252	11798	0.398	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.70	276	13426	0.395	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.73	278	11170	0.394	ng/ul#	95
26) Benzo(g,h,i)perylene	27.53	276	11174	0.391	ng/ul	96

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

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