

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080817\
 Data File : BE093724.D
 Acq On : 8 Aug 2017 19:17
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.445

Manual Integrations
 APPROVED

Sohil
 8/9/2017 7:14:38 PM

Quant Time: Aug 09 05:28:42 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 09 04:20:36 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	13863	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	7833	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	18937	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19685	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	17355m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8817	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	20803	0.35	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	13794	0.409	ng/ul#	92
5) 2-Methylnaphthalene	12.36	142	9617	0.391	ng/ul	98
7) Acenaphthylene	14.24	152	13307	0.406	ng/ul	97
8) Acenaphthene	14.59	153	10363	0.402	ng/ul	98
9) Fluorene	15.57	166	12032	0.369	ng/ul#	90
11) Pentachlorophenol	16.91	266	850	0.272	ng/ul	97
12) Phenanthrene	17.29	178	20280	0.393	ng/ul#	90
13) Anthracene	17.38	178	18640	0.371	ng/ul#	89
15) Fluoranthene	19.30	202	22402	0.347	ng/ul	84
17) Pyrene	19.66	202	22885	0.398	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	21441	0.376	ng/ul#	85
19) Chrysene	21.45	228	21844	0.399	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	20670	0.385	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	21475m	0.406	ng/ul	
23) Benzo(a)pyrene	23.80	252	19859m	0.389	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	22110m	0.392	ng/ul	
25) Dibenzo(a,h)anthracene	26.57	278	18371m	0.390	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	17955m	0.378	ng/ul	

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

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