

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094655.D
 Acq On : 31 Oct 2017 11:26
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.468

Manual Integrations
 APPROVED

Sohil
 11/1/2017 9:31:48 AM

Quant Time: Nov 01 01:35:17 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 04:29:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1223	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7515	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4192	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8850	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9028	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8712	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4860	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	10628	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	7121	0.388	ng/ul#	95
5) 2-Methylnaphthalene	12.35	142	4891	0.380	ng/ul	97
7) Acenaphthylene	14.24	152	7602	0.401	ng/ul	99
8) Acenaphthene	14.57	153	5669	0.416	ng/ul	96
9) Fluorene	15.57	166	6643	0.434	ng/ul	92
11) Pentachlorophenol	17.00	266	179	0.405	ng/ul	100
12) Phenanthrene	17.29	178	11814m	0.505	ng/ul	
13) Anthracene	17.39	178	9596	0.416	ng/ul	96
15) Fluoranthene	19.31	202	10907	0.404	ng/ul	84
17) Pyrene	19.68	202	12066	0.432	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	10599	0.411	ng/ul#	87
19) Chrysene	21.47	228	10258	0.385	ng/ul	99
21) Benzo(b)fluoranthene	23.19	252	9777	0.399	ng/ul	92
22) Benzo(k)fluoranthene	23.23	252	10386	0.395	ng/ul	94
23) Benzo(a)pyrene	23.86	252	9695	0.396	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.67	276	9050	0.357	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.67	278	7979	0.374	ng/ul	94
26) Benzo(g,h,i)perylene	27.52	276	6644	0.326	ng/ul	93

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

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