

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE100717\
 Data File : BE094204.D
 Acq On : 7 Oct 2017 16:03
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.419

Manual Integrations
 APPROVED

Sohil

Quant Time: Oct 08 01:33:41 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 01:38:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	10/9/2017 6:52:55 PM
1) 1,4-Dichlorobenzene-d4	7.90	152	1018	0.40	ng/ul	0.00	
2) Naphthalene-d8	10.69	136	5257	0.40	ng/ul	0.00	
6) Acenaphthene-d10	14.51	164	3262	0.40	ng/ul	-0.02	
10) Phenanthrene-d10	17.24	188	7830	0.40	ng/ul	-0.03	
16) Chrysene-d12	21.40	240	7514	0.40	ng/ul	-0.01	
20) Perylene-d12	23.91	264	8090	0.40	ng/ul	-0.02	
System Monitoring Compounds							
4) 2-Methylnaphthalene-d10	12.28	152	3382	0.41	ng/ul	-0.01	
14) Fluoranthene-d10	19.26	212	9351	0.36	ng/ul	-0.02	
Target Compounds				Qvalue			
3) Naphthalene	10.74	128	5028	0.39	ng/ul#	93	
5) 2-Methylnaphthalene	12.35	142	3585	0.43	ng/ul	99	
7) Acenaphthylene	14.24	152	6012	0.47	ng/ul	97	
8) Acenaphthene	14.57	153	4252	0.41	ng/ul	97	
9) Fluorene	15.56	166	5123	0.42	ng/ul	92	
11) Pentachlorophenol	16.93	266	468	0.51	ng/ul	92	
12) Phenanthrene	17.27	178	7967m	0.39	ng/ul		
13) Anthracene	17.38	178	8156	0.41	ng/ul	99	
15) Fluoranthene	19.30	202	9432	0.35	ng/ul	87	
17) Pyrene	19.66	202	10236	0.49	ng/ul#	84	
18) Benzo(a)anthracene	21.39	228	8819	0.44	ng/ul#	80	
19) Chrysene	21.44	228	8163	0.35	ng/ul#	80	
21) Benzo(b)fluoranthene	23.14	252	8476	0.39	ng/ul#	1	
22) Benzo(k)fluoranthene	23.19	252	9221	0.36	ng/ul#	1	
23) Benzo(a)pyrene	23.80	252	8267	0.36	ng/ul#	1	
24) Indeno(1,2,3-cd)pyrene	26.55	276	9059	0.39	ng/ul#	81	
25) Dibenzo(a,h)anthracene	26.57	278	7549	0.39	ng/ul#	1	
26) Benzo(g,h,i)perylene	27.37	276	8377	0.42	ng/ul#	28	

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

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