

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094380.D
 Acq On : 15 Oct 2017 6:25
 Operator : SJ/JU
 Sample : I5548-16
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AF6

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:47 PM

Quant Time: Oct 16 06:25:21 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1608	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	8586	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5239	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11977	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	11853	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	11140m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2929	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	7609	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	1802	0.086	ng/ul#	89
5) 2-Methylnaphthalene	12.35	142	1402	0.103	ng/ul	98
7) Acenaphthylene	14.24	152	3256	0.159	ng/ul	96
8) Acenaphthene	14.58	153	2147	0.128	ng/ul	98
9) Fluorene	15.57	166	2758	0.140	ng/ul	93
12) Phenanthrene	17.29	178	56559	1.807	ng/ul#	88
13) Anthracene	17.38	178	13083	0.430	ng/ul#	89
15) Fluoranthene	19.31	202	140804	3.424	ng/ul	84
17) Pyrene	19.67	202	135340	4.104	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	72614	2.309	ng/ul#	88
19) Chrysene	21.46	228	73663	2.029	ng/ul	97
21) Benzo(b)fluoranthene	23.17	252	91987m	3.040	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	34742m	0.993	ng/ul	
23) Benzo(a)pyrene	23.84	252	62951m	2.002	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.63	276	41996m	1.306	ng/ul	
25) Dibenzo(a,h)anthracene	26.63	278	11927m	0.442	ng/ul	
26) Benzo(a,h,i)perylene	27.46	276	33987m	1.225	ng/ul	

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

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