

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094305.D
 Acq On : 12 Oct 2017 6:56
 Operator : SJ/JU
 Sample : I5550-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AF1

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:06:14 PM

Quant Time: Oct 12 09:41:59 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 08:55:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1773	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	9094	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	5257	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	10796m	0.40	ng/ul	0.02
16) Chrysene-d12	21.44	240	9687	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11953m	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4809	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	11534	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	17223	0.773	ng/ul#	91
5) 2-Methylnaphthalene	12.37	142	6875	0.476	ng/ul	98
7) Acenaphthylene	14.26	152	2299	0.112	ng/ul#	81
8) Acenaphthene	14.60	153	17188	1.019	ng/ul	99
9) Fluorene	15.58	166	25999	1.316	ng/ul	92
12) Phenanthrene	17.31	178	425160	15.070	ng/ul#	89
13) Anthracene	17.40	178	104981	3.826	ng/ul#	88
15) Fluoranthene	19.33	202	479476	12.935	ng/ul	84
17) Pyrene	19.69	202	408274	15.148	ng/ul#	84
18) Benzo(a)anthracene	21.42	228	209108	8.137	ng/ul#	86
19) Chrysene	21.47	228	178180	6.004	ng/ul	96
21) Benzo(b)fluoranthene	23.19	252	166340	5.123	ng/ul	91
22) Benzo(k)fluoranthene	23.23	252	65731m	1.751	ng/ul	
23) Benzo(a)pyrene	23.86	252	129336m	3.833	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.66	276	56611m	1.640	ng/ul	
25) Dibenzo(a,h)anthracene	26.67	278	14844m	0.513	ng/ul	
26) Benzo(a,h,i)perylene	27.50	276	38592m	1.296	ng/ul	

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

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