

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094308.D
 Acq On : 12 Oct 2017 8:47
 Operator : SJ/JU
 Sample : I5649-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-39(5-7)-100317

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:06:20 PM

Quant Time: Oct 12 20:20:58 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	3616	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	11002	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5742	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	11536m	0.40	ng/ul	0.02
16) Chrysene-d12	21.43	240	10562	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8599m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6648	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	13706	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	20404	0.757	ng/ul#	63
5) 2-Methylnaphthalene	12.36	142	1002	0.057	ng/ul	98
7) Acenaphthylene	14.26	152	2158	0.096	ng/ul#	45
8) Acenaphthene	14.60	153	4258	0.231	ng/ul	97
9) Fluorene	15.58	166	3813	0.177	ng/ul	92
12) Phenanthrene	17.31	178	18513	0.614	ng/ul	95
13) Anthracene	17.40	178	6435m	0.219	ng/ul	
15) Fluoranthene	19.32	202	32376	0.817	ng/ul	84
17) Pyrene	19.69	202	37650	1.281	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	16352	0.584	ng/ul#	83
19) Chrysene	21.47	228	15890	0.491	ng/ul#	78
21) Benzo(b)fluoranthene	23.19	252	18227	0.780	ng/ul#	73
22) Benzo(k)fluoranthene	23.23	252	7414m	0.275	ng/ul	
23) Benzo(a)pyrene	23.85	252	16082m	0.663	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.65	276	8469m	0.341	ng/ul	
26) Benzo(a,h,i)perylene	27.49	276	7831m	0.366	ng/ul	

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

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