

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101117\
 Data File : BE094307.D
 Acq On : 12 Oct 2017 8:11
 Operator : SJ/JU
 Sample : I5649-09
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-57(2-4)-100317

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:06:17 PM

Quant Time: Oct 12 20:16:25 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	1690	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	8710	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	5028	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	10508m	0.40	ng/ul	0.02
16) Chrysene-d12	21.43	240	9822	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8555m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4307	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9719	0.28	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	9104	0.427	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	2080	0.150	ng/ul	98
7) Acenaphthylene	14.26	152	890	0.045	ng/ul#	73
8) Acenaphthene	14.60	153	613	0.038	ng/ul#	91
9) Fluorene	15.58	166	1906	0.101	ng/ul#	88
12) Phenanthrene	17.31	178	22061	0.803	ng/ul#	93
13) Anthracene	17.39	178	2699	0.101	ng/ul#	91
15) Fluoranthene	19.32	202	25389	0.704	ng/ul	84
17) Pyrene	19.68	202	27519	1.007	ng/ul#	84
18) Benzo(a)anthracene	21.42	228	15426	0.592	ng/ul#	68
19) Chrysene	21.47	228	25728	0.855	ng/ul#	78
21) Benzo(b)fluoranthene	23.19	252	24258m	1.044	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	7248m	0.270	ng/ul	
23) Benzo(a)pyrene	23.86	252	16462m	0.682	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.65	276	7950m	0.322	ng/ul	
26) Benzo(a,h,i)perylene	27.49	276	8730m	0.410	ng/ul	

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

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