

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012898.D
 Acq On : 11 Nov 2017 12:53
 Operator : SJ/JU
 Sample : I6317-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHJO

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:51:04 PM

Quant Time: Nov 13 01:24:19 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 01:07:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	57094	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	219411	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	132919	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	312439	20.00	ng/ul	0.00
17) Chrysene-d12	21.35	240	392704	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	430232	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	348732	25.22	ng/ul	0.00
10) Fluorene-d10	15.42	176	257615	27.24	ng/ul	0.00
14) Anthracene-d10	17.27	188	399964	26.48	ng/ul	0.00
18) Pyrene-d10	19.56	212	457938	26.06	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	537539	26.46	ng/ul	0.00
Target Compounds						
16) Fluoranthene	19.23	202	46777	2.203	ng/ul#	90
19) Pyrene	19.59	202	43295	1.921	ng/ul#	93
20) Benzo(a)anthracene	21.34	228	31147m	1.289	ng/ul	
21) Chrysene	21.39	228	34672m	1.606	ng/ul	
23) Benzo(b)fluoranthene	22.94	252	68667	2.684	ng/ul#	97
26) Benzo(a)pyrene	23.53	252	37440	1.509	ng/ul	97

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

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