

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012919.D
 Acq On : 12 Nov 2017 01:19
 Operator : SJ/JU
 Sample : I6322-03MSD
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHH9MSD

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:49:53 PM

Quant Time: Nov 13 03:20:19 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 02:06:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	84862	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	317572	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	185345	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	435395	20.00	ng/ul	0.00
17) Chrysene-d12	21.35	240	534692	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	603026	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	483677	25.09	ng/ul	0.00
10) Fluorene-d10	15.42	176	353236	26.79	ng/ul	0.00
14) Anthracene-d10	17.27	188	548941	26.08	ng/ul	0.00
18) Pyrene-d10	19.56	212	610123	25.50	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	739776	25.98	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthene	14.49	153	300619	23.930	ng/ul	96
13) Phenanthrene	17.21	178	334130	13.894	ng/ul	99
15) Anthracene	17.30	178	73565	2.944	ng/ul	97
16) Fluoranthene	19.23	202	791378	26.751	ng/ul#	92
19) Pyrene	19.59	202	1236764	40.295	ng/ul#	94
20) Benzo(a)anthracene	21.34	228	439648	13.363	ng/ul	98
21) Chrysene	21.39	228	385280	13.104	ng/ul	96
23) Benzo(b)fluoranthene	22.94	252	672157m	18.744	ng/ul	
24) Benzo(k)fluoranthene	22.99	252	209878m	6.124	ng/ul	
26) Benzo(a)pyrene	23.53	252	404648m	11.635	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.93	276	338800	7.553	ng/ul	95
28) Dibenzo(a,h)anthracene	25.93	278	98345	2.608	ng/ul#	76
29) Benzo(a,h,i)perylene	26.63	276	256944	6.862	ng/ul#	94

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

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