

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031518\
 Data File : BM014589.D
 Acq On : 15 Mar 2018 22:51
 Operator : SJ/JU
 Sample : J1904-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEL27

Quant Time: Mar 26 12:08:18 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031518-PAH.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 16 15:09:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	103503	20.00	ng/ul	0.00
2) Naphthalene-d8	10.22	136	451165	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	265810	20.00	ng/ul	0.00
12) Phenanthrene-d10	16.89	188	549565	20.00	ng/ul	0.00
18) Chrysene-d12	21.11	240	433618	20.00	ng/ul	0.00
23) Perylene-d12	23.27	264	374750	20.00	ng/ul	-0.01
System Monitoring Compounds						
3) 2,4-Dichlorophenol-d3	9.87	165	220028	31.56	ng/ul	-0.01
7) Acenaphthylene-d8	13.81	160	955941	33.65	ng/ul	0.00
10) Fluorene-d10	15.13	176	626186	33.65	ng/ul	0.00
15) Anthracene-d10	16.99	188	887592	33.74	ng/ul	0.00
19) Pyrene-d10	19.30	212	902435	34.62	ng/ul	0.00
26) Benzo(a)pyrene-d12	23.13	264	611289	33.72	ng/ul	0.00
Target Compounds				Qvalue		
14) Phenanthrene	16.93	178	52754	1.714	ng/ul	99
17) Fluoranthene	18.97	202	73992	2.118	ng/ul#	97
20) Pyrene	19.33	202	84906	2.604	ng/ul#	95
22) Chrysene	21.14	228	43433	1.723	ng/ul	93
24) Benzo(b)fluoranthene	22.63	252	38395	1.630	ng/ul#	86

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

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