

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091219\
 Data File : BM022665.D
 Acq On : 12 Sep 2019 13:44
 Operator : HP/JU
 Sample : K4706-08DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BF574DL

Quant Time: Sep 12 15:00:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	10309	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	42196	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	22831	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	53714	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	43925	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	41612	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	3860	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	9772	0.07	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.69	128	2773	0.023	ng/ul#	89
5) 2-Methylnaphthalene	12.30	142	7324	0.089	ng/ul	100
8) Acenaphthene	14.54	153	25457	0.288	ng/ul	100
9) Fluorene	15.52	166	39154	0.386	ng/ul	100
12) Phenanthrene	17.25	178	131277	0.767	ng/ul	100
13) Anthracene	17.34	178	7915	0.054	ng/ul	95
15) Fluoranthene	19.27	202	26190	0.142	ng/ul	99
17) Pyrene	19.63	202	21770	0.109	ng/ul	98
18) Benzo(a)anthracene	21.38	228	3846	0.023	ng/ul	95
19) Chrysene	21.38	228	3846	0.022	ng/ul	97
21) Benzo(b)fluoranthene	22.99	252	6333	0.035	ng/ul#	18
22) Benzo(k)fluoranthene	22.99	252	6258	0.034	ng/ul#	17

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

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