

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025270.D
 Acq On : 05 Mar 2020 03:12
 Operator : CG/JU
 Sample : L1516-04DL 5X
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD22DL

Quant Time: Mar 05 04:28:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 05 00:55:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3664	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	15422	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11268	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	27739	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	23953	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	19883	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1349	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	3775	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.02	152	1127	0.021	ng/ul#	93
8) Acenaphthene	14.36	153	25472	0.622	ng/ul	99
9) Fluorene	15.35	166	18922	0.369	ng/ul	97
12) Phenanthrene	17.08	178	74384	0.832	ng/ul	95
13) Anthracene	17.17	178	8166	0.103	ng/ul	98
15) Fluoranthene	19.10	202	114513	1.023	ng/ul	94
17) Pyrene	19.46	202	74114	0.816	ng/ul	95
18) Benzo(a)anthracene	21.21	228	10195	0.116	ng/ul	99
19) Chrysene	21.26	228	9237	0.095	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	3851	0.048	ng/ul	72
22) Benzo(k)fluoranthene	22.81	252	1905	0.023	ng/ul#	27
23) Benzo(a)pyrene	23.32	252	1592	0.024	ng/ul#	44

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

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