

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025161.D
 Acq On : 29 Feb 2020 10:54
 Operator : CG/JU
 Sample : L1542-03
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD39

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:34 AM

Quant Time: Mar 02 01:43:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 01:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2587	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10094	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6628	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14731m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	15604	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17827	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7738	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	19901	0.39	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.03	152	908	0.024	ng/ul#	71
9) Fluorene	15.37	166	781m	0.024	ng/ul	
12) Phenanthrene	17.10	178	2176	0.042	ng/ul#	91
15) Fluoranthene	19.11	202	3060	0.041	ng/ul	86
17) Pyrene	19.47	202	3391	0.048	ng/ul#	86
18) Benzo(a)anthracene	21.22	228	1418	0.021	ng/ul#	82
19) Chrysene	21.27	228	1948	0.029	ng/ul#	89
21) Benzo(b)fluoranthene	22.78	252	1761m	0.025	ng/ul	

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

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