

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025155.D
 Acq On : 29 Feb 2020 04:56
 Operator : CG/JU
 Sample : L1506-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCY5

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:28 AM

Quant Time: Mar 02 00:56:54 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 00:36:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2798	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10989	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7146	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16495	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	16092	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17399	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	3766	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	10708	0.19	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.03	152	1447	0.035	ng/ul#	92
9) Fluorene	15.37	166	1922	0.055	ng/ul	96
12) Phenanthrene	17.10	178	128811	2.220	ng/ul	95
13) Anthracene	17.19	178	63264	1.125	ng/ul	95
15) Fluoranthene	19.11	202	424993	5.043	ng/ul	96
17) Pyrene	19.47	202	335287	4.624	ng/ul	95
18) Benzo(a)anthracene	21.22	228	157452	2.266	ng/ul	98
19) Chrysene	21.27	228	140037	2.024	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	121994m	1.789	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	43706m	0.647	ng/ul	
23) Benzo(a)pyrene	23.34	252	68290	1.097	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.62	276	27733	0.355	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	7825	0.119	ng/ul#	85
26) Benzo(g,h,i)perylene	26.28	276	17231	0.262	ng/ul#	93

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

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