

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025140.D
 Acq On : 28 Feb 2020 19:21
 Operator : CG/JU
 Sample : L1507-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD04

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:38:38 AM

Quant Time: Feb 29 05:28:37 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 : Sat Feb 29 04:43:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2650	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10377	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6971	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16068m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17333	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	19539	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6606	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	20563	0.37	ng/ul	0.00
Target Compounds				Ovalue		
15) Fluoranthene	19.11	202	9016	0.110	ng/ul	98
17) Pyrene	19.47	202	9898	0.127	ng/ul	98
18) Benzo(a)anthracene	21.22	228	3988	0.053	ng/ul	97
19) Chrysene	21.27	228	7991	0.107	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	8050m	0.105	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	3045m	0.040	ng/ul	
23) Benzo(a)pyrene	23.34	252	2129	0.030	ng/ul#	67
24) Indeno(1,2,3-cd)pyrene	25.62	276	1918	0.022	ng/ul#	100
26) Benzo(g,h,i)perylene	26.29	276	1612	0.022	ng/ul#	81

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

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