

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040320\
 Data File : BM025898.D
 Acq On : 02 Apr 2020 12:06
 Operator : CG/JU
 Sample : L1516-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD24

Manual Integrations
 APPROVED

mohammad
 4/3/2020 8:05:11 AM

Quant Time: Apr 02 12:45:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2405	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	10226	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7312	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	17259m	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	19572	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	19860	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	4860	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	13897	0.26	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.99	178	166925	3.070	ng/ul	98
15) Fluoranthene	19.01	202	432186	6.433	ng/ul	99
17) Pyrene	19.38	202	356081	5.080	ng/ul	99
18) Benzo(a)anthracene	21.13	228	48434	0.692	ng/ul	100
19) Chrysene	21.18	228	77382	1.013	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	22492m	0.308	ng/ul	
22) Benzo(k)fluoranthene	22.70	252	10793m	0.138	ng/ul	
23) Benzo(a)pyrene	23.20	252	10630	0.163	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.41	276	13654	0.156	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.42	278	3987	0.055	ng/ul#	82
26) Benzo(g,h,i)perylene	26.06	276	11014	0.149	ng/ul	96

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

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