

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040320\
 Data File : BM025900.D
 Acq On : 02 Apr 2020 13:53
 Operator : CG/JU
 Sample : L1516-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD23

Manual Integrations
 APPROVED

mohammad
 4/3/2020 8:05:14 AM

Quant Time: Apr 02 14:36:12 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	7.56	152	2336	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9983	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7055	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	16647m	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18441	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	19967	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	4318	0.24	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	12396	0.24	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.26	153	652	0.026	ng/ul#	90
9) Fluorene	15.26	166	666	0.021	ng/ul#	97
12) Phenanthrene	16.99	178	181852	3.468	ng/ul	98
15) Fluoranthene	19.01	202	341207	5.265	ng/ul	99
17) Pyrene	19.37	202	329520	4.989	ng/ul	99
18) Benzo(a)anthracene	21.13	228	4165	0.063	ng/ul	96
19) Chrysene	21.18	228	4864	0.068	ng/ul	95
21) Benzo(b)fluoranthene	22.66	252	8781m	0.119	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	3497m	0.044	ng/ul	
23) Benzo(a)pyrene	23.20	252	3850	0.059	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	25.42	276	2424	0.028	ng/ul#	63
26) Benzo(g,h,i)perylene	26.06	276	1931	0.026	ng/ul#	75

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

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