

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025241.D
 Acq On : 04 Mar 2020 09:16
 Operator : CG/JU
 Sample : L1516-06
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD24

Manual Integrations
 APPROVED

mohammad
 3/5/2020 7:56:56 AM

Quant Time: Mar 04 11:37:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	4660	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	20391	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	15143	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	36686m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	35126	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	32374	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	8338	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	24385	0.20	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	5454	0.046	ng/ul	99
15) Fluoranthene	19.10	202	15361	0.104	ng/ul	97
17) Pyrene	19.47	202	14777m	0.111	ng/ul	
18) Benzo(a)anthracene	21.22	228	9140	0.071	ng/ul	97
19) Chrysene	21.27	228	23587	0.166	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	32178	0.248	ng/ul	95
22) Benzo(k)fluoranthene	22.81	252	12137	0.089	ng/ul#	80
23) Benzo(a)pyrene	23.33	252	14429	0.133	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.60	276	16060	0.118	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	4583	0.041	ng/ul#	85
26) Benzo(g,h,i)perylene	26.27	276	12419	0.108	ng/ul#	94

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

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