

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025214.D
 Acq On : 03 Mar 2020 16:21
 Operator : CG/JU
 Sample : L1507-04DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD02DL

Manual Integrations
 APPROVED

mohammad
 3/4/2020 7:52:47 AM

Quant Time: Mar 03 17:07:22 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 : Tue Mar 03 14:30:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	3409	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	13363	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	9261	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	22205	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	22243	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	20646	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	863	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2534	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	5764	0.080	ng/ul	95
15) Fluoranthene	19.11	202	15076	0.168	ng/ul	95
17) Pyrene	19.47	202	29492	0.350	ng/ul#	94
18) Benzo(a)anthracene	21.22	228	19117	0.234	ng/ul	99
19) Chrysene	21.27	228	14540	0.162	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	34138m	0.413	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	15990m	0.185	ng/ul	
23) Benzo(a)pyrene	23.34	252	16116	0.232	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.62	276	9587	0.111	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	3598	0.050	ng/ul#	87
26) Benzo(g,h,i)perylene	26.28	276	5960	0.081	ng/ul#	93

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

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