

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027972.D
 Acq On : 25 Nov 2020 20:13
 Operator : JU/CG
 Sample : L4749-06DL 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B69DL

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:43:59 AM

Quant Time: Nov 26 01:06:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 26 00:48:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	1096	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4427	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2496	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4584	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2493	0.40	ng/ul	0.00
20) Perylene-d12	24.17	264	2346	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	90	0.01	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	140	0.01	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.69	178	4111	0.284	ng/ul	100
13) Anthracene	17.78	178	642	0.046	ng/ul#	79
15) Fluoranthene	19.67	202	12565	0.783	ng/ul	97
17) Pyrene	20.03	202	10042	0.793	ng/ul#	95
18) Benzo(a)anthracene	21.74	228	4554	0.453	ng/ul	95
19) Chrysene	21.80	228	5657	0.577	ng/ul	96
21) Benzo(b)fluoranthene	23.44	252	7002m	0.710	ng/ul	
22) Benzo(k)fluoranthene	23.48	252	2635m	0.273	ng/ul	
23) Benzo(a)pyrene	24.07	252	4387	0.509	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	26.66	276	3293	0.353	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.67	278	828	0.107	ng/ul#	40
26) Benzo(g,h,i)perylene	27.44	276	2958	0.378	ng/ul#	92

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

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