

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027957.D
 Acq On : 25 Nov 2020 11:05
 Operator : JU/CG
 Sample : L4751-04DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B84DL

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:43:03 AM

Quant Time: Nov 25 12:01:36 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 : Wed Nov 25 10:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	1242	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4808	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2496	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4107	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2699	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2790	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	176	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	240	0.02	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	15.00	153	470	0.052	ng/ul	96
9) Fluorene	15.98	166	912	0.086	ng/ul#	97
12) Phenanthrene	17.70	178	22652	1.746	ng/ul	97
13) Anthracene	17.79	178	6533	0.522	ng/ul	100
15) Fluoranthene	19.68	202	76345	5.307	ng/ul	100
17) Pyrene	20.03	202	55407	4.042	ng/ul	98
18) Benzo(a)anthracene	21.74	228	31776	2.918	ng/ul	100
19) Chrysene	21.80	228	34672	3.269	ng/ul	98
21) Benzo(b)fluoranthene	23.44	252	47992m	4.093	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	16655m	1.450	ng/ul	
23) Benzo(a)pyrene	24.08	252	25562	2.493	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.67	276	19372	1.745	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.68	278	5087	0.550	ng/ul	99
26) Benzo(g,h,i)perylene	27.45	276	15771	1.693	ng/ul	95

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

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