

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027942.D
 Acq On : 24 Nov 2020 14:43
 Operator : JU/CG
 Sample : L4749-12
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AX9

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:46:04 PM

Quant Time: Nov 24 15:16:23 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 : Tue Nov 24 15:00:24 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1006	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	3947	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2179	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3596	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2013	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	1907	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	931	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	1301	0.14	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	303	0.026	ng/ul#	54
5) 2-Methylnaphthalene	12.79	142	198	0.025	ng/ul	95
12) Phenanthrene	17.70	178	3738	0.329	ng/ul	99
13) Anthracene	17.79	178	473	0.043	ng/ul#	71
15) Fluoranthene	19.68	202	9345	0.742	ng/ul	97
17) Pyrene	20.03	202	7654	0.749	ng/ul	96
18) Benzo(a)anthracene	21.74	228	3359	0.414	ng/ul	94
19) Chrysene	21.80	228	4099	0.518	ng/ul	96
21) Benzo(b)fluoranthene	23.44	252	5577m	0.696	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	2056m	0.262	ng/ul	
23) Benzo(a)pyrene	24.07	252	3449	0.492	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	26.67	276	2540	0.335	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.68	278	628	0.099	ng/ul#	17
26) Benzo(g,h,i)perylene	27.45	276	2292	0.360	ng/ul#	89

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

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