

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027948.D
 Acq On : 24 Nov 2020 18:20
 Operator : JU/CG
 Sample : L4749-16
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B15

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:46:20 PM

Quant Time: Nov 25 01:13:50 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	864	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	3515	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2081	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3522m	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2030	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	1884	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1694	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	2322	0.25	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.70	178	1618	0.145	ng/ul	94
13) Anthracene	17.79	178	362	0.034	ng/ul#	68
15) Fluoranthene	19.68	202	2151	0.174	ng/ul	89
17) Pyrene	20.03	202	2093	0.203	ng/ul#	85
18) Benzo(a)anthracene	21.75	228	733	0.089	ng/ul#	79
19) Chrysene	21.80	228	741	0.093	ng/ul#	77
21) Benzo(b)fluoranthene	23.45	252	770m	0.097	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	325m	0.042	ng/ul	
23) Benzo(a)pyrene	24.08	252	553	0.080	ng/ul#	27
24) Indeno(1,2,3-cd)pyrene	26.67	276	364	0.049	ng/ul#	65
26) Benzo(g,h,i)perylene	27.45	276	326	0.052	ng/ul#	30

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

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