

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027921.D
 Acq On : 23 Nov 2020 19:49
 Operator : JU/CG
 Sample : L4711-08DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B33DL

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:44:55 PM

Quant Time: Nov 24 01:04:34 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 00:42:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1197	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4708	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2660	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4889	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	3010	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2643	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	264	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	434	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	1229	0.087	ng/ul#	90
5) 2-Methylnaphthalene	12.80	142	782	0.081	ng/ul	99
7) Acenaphthylene	14.67	152	343	0.026	ng/ul#	61
8) Acenaphthene	15.01	153	4847	0.498	ng/ul	100
9) Fluorene	15.98	166	3987	0.353	ng/ul	99
12) Phenanthrene	17.70	178	71487	4.629	ng/ul	97
13) Anthracene	17.79	178	13948	0.936	ng/ul	99
15) Fluoranthene	19.68	202	125070	7.304	ng/ul	99
17) Pyrene	20.03	202	101824	6.661	ng/ul	98
18) Benzo(a)anthracene	21.75	228	43219	3.559	ng/ul	100
19) Chrysene	21.80	228	46948	3.969	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	50667m	4.561	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	17832m	1.639	ng/ul	
23) Benzo(a)pyrene	24.08	252	33486	3.448	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.67	276	21795	2.073	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.68	278	5523	0.631	ng/ul	96
26) Benzo(a,h,i)perylene	27.45	276	19208	2.177	ng/ul	93

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

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