

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027902.D
 Acq On : 22 Nov 2020 05:01
 Operator : JU/CG
 Sample : L4711-14
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B66

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:59:35 AM

Quant Time: Nov 23 06:05:18 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 : Mon Nov 23 05:39:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1106	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4362	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2638	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4440m	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2349	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2093	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2269	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3367	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	560	0.043	ng/ul#	76
5) 2-Methylnaphthalene	12.80	142	373	0.042	ng/ul	98
7) Acenaphthylene	14.67	152	435	0.034	ng/ul#	65
8) Acenaphthene	15.01	153	458	0.047	ng/ul#	92
9) Fluorene	15.98	166	472	0.042	ng/ul#	96
12) Phenanthrene	17.70	178	15729	1.121	ng/ul	98
13) Anthracene	17.79	178	2156	0.159	ng/ul	94
15) Fluoranthene	19.68	202	38468	2.474	ng/ul	100
17) Pyrene	20.04	202	30656	2.570	ng/ul	99
18) Benzo(a)anthracene	21.74	228	12450	1.314	ng/ul	99
19) Chrysene	21.80	228	16207	1.756	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	19066m	2.168	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	6825m	0.792	ng/ul	
23) Benzo(a)pyrene	24.07	252	11991	1.559	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.68	276	8906	1.070	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.68	278	2214	0.319	ng/ul#	67
26) Benzo(a,h,i)perylene	27.45	276	8599	1.231	ng/ul	98

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

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