

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027980.D
 Acq On : 26 Nov 2020 01:03
 Operator : JU/CG
 Sample : L4749-19DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B18DL

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:44:24 AM

Quant Time: Nov 26 01:51:40 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 : Thu Nov 26 00:48:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	1412	0.40	ng/ul	0.00
2) Naphthalene-d8	11.15	136	5727	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	3148	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	5489	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	3150	0.40	ng/ul	0.00
20) Perylene-d12	24.17	264	2957	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	195	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	298	0.02	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.66	152	356	0.023	ng/ul#	65
8) Acenaphthene	15.00	153	472	0.041	ng/ul	96
9) Fluorene	15.97	166	557	0.042	ng/ul#	95
11) Pentachlorophenol	17.30	266	43	0.023	ng/ul	98
12) Phenanthrene	17.69	178	12706	0.733	ng/ul	98
13) Anthracene	17.78	178	1895	0.113	ng/ul	94
15) Fluoranthene	19.67	202	22132	1.151	ng/ul	98
17) Pyrene	20.03	202	18292	1.143	ng/ul	97
18) Benzo(a)anthracene	21.74	228	7723	0.608	ng/ul	98
19) Chrysene	21.79	228	8946	0.723	ng/ul	97
21) Benzo(b)fluoranthene	23.44	252	10109m	0.813	ng/ul	
22) Benzo(k)fluoranthene	23.48	252	3568m	0.293	ng/ul	
23) Benzo(a)pyrene	24.06	252	6833	0.629	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	26.66	276	4433	0.377	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.66	278	1172	0.120	ng/ul#	63
26) Benzo(a,h,i)perylene	27.43	276	3973	0.403	ng/ul	97

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

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