

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082020\
 Data File : BM027159.D
 Acq On : 20 Aug 2020 16:23
 Operator : JU/CG
 Sample : L1955-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-SOIL-07

Quant Time: Aug 20 16:57:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 06:55:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	511	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	1642	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	1062	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	2458	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	2382	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	2497	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	974	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	2752	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	297	0.060	ng/ul#	47
5) 2-Methylnaphthalene	12.08	142	193	0.057	ng/ul	95
7) Acenaphthylene	13.98	152	275	0.060	ng/ul#	65
8) Acenaphthene	14.32	153	251	0.063	ng/ul	90
9) Fluorene	15.32	166	290	0.060	ng/ul#	90
11) Pentachlorophenol	16.69	266	51	0.063	ng/ul#	70
12) Phenanthrene	17.05	178	522	0.063	ng/ul#	70
13) Anthracene	17.14	178	403	0.053	ng/ul#	59
15) Fluoranthene	19.07	202	630	0.063	ng/ul	74
17) Pyrene	19.43	202	634	0.062	ng/ul#	78
18) Benzo(a)anthracene	21.19	228	569	0.060	ng/ul#	82
19) Chrysene	21.24	228	736	0.072	ng/ul#	90
21) Benzo(b)fluoranthene	22.74	252	719	0.067	ng/ul#	44
22) Benzo(k)fluoranthene	22.79	252	782	0.070	ng/ul#	47
23) Benzo(a)pyrene	23.30	252	608	0.063	ng/ul#	38
24) Indeno(1,2,3-cd)pyrene	25.58	276	822	0.061	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.60	278	639	0.059	ng/ul#	29
26) Benzo(g,h,i)perylene	26.24	276	717	0.062	ng/ul#	70

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

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