

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM123020\
 Data File : BM028286.D
 Acq On : 30 Dec 2020 15:13
 Operator : JU/CG
 Sample : L5175-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2W00

Quant Time: Dec 30 15:56:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM121120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 30 10:02:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	1710	0.40	ng/ul	0.00
4) Naphthalene-d8	11.01	136	6653	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.81	164	3417	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.53	188	6817	0.40	ng/ul	0.00
17) Chrysene-d12	21.65	240	6059	0.40	ng/ul	0.00
23) Perylene-d12	24.00	264	5894	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.41	96	780	0.40	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.58	152	2546	0.25	ng/ul	0.00
18) Fluoranthene-d10	19.53	212	5188	0.29	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.06	128	8653	0.461	ng/ul	99
7) 2-Methylnaphthalene	12.65	142	6202	0.501	ng/ul	99
10) Acenaphthylene	14.53	152	7407	0.537	ng/ul	99
11) Acenaphthene	14.87	153	7478	0.617	ng/ul	100
14) Pentachlorophenol	17.18	266	1344	0.777	ng/ul	99
16) Anthracene	17.66	178	9096	0.496	ng/ul	99
20) Pyrene	19.92	202	17176	0.722	ng/ul	96
22) Chrysene	21.68	228	13217	0.614	ng/ul	100
24) Benzo(b)fluoranthene	23.29	252	16095	0.679	ng/ul	99
25) Benzo(k)fluoranthene	23.33	252	25020	1.098	ng/ul	95
26) Benzo(a)pyrene	23.90	252	8927	0.456	ng/ul	96
28) Dibenzo(a,h)anthracene	26.43	278	15833	0.795	ng/ul	97

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

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