

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121420\
 Data File : BM028116.D
 Acq On : 14 Dec 2020 17:42
 Operator : JU/CG
 Sample : L4958-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2S43

Quant Time: Dec 14 18:16:40 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 : Mon Dec 14 12:18:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	2463	0.40	ng/ul	0.00
4) Naphthalene-d8	11.05	136	10177	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.84	164	5887	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.56	188	12559	0.40	ng/ul	0.00
17) Chrysene-d12	21.67	240	10408	0.40	ng/ul	0.00
23) Perylene-d12	24.04	264	8365	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.43	96	1183	0.42	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.62	152	4452	0.29	ng/ul	0.00
18) Fluoranthene-d10	19.56	212	9887	0.32	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.10	128	16595	0.578	ng/ul	100
7) 2-Methylnaphthalene	12.69	142	11211	0.592	ng/ul	99
12) Fluorene	15.88	166	16154	0.680	ng/ul	99
14) Pentachlorophenol	17.21	266	5336	1.674	ng/ul	99
15) Phenanthrene	17.60	178	27249	0.707	ng/ul	100
19) Fluoranthene	19.59	202	25829	0.631	ng/ul	99
21) Benzo(a)anthracene	21.66	228	28233	0.823	ng/ul	99
25) Benzo(k)fluoranthene	23.37	252	23333	0.722	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.47	276	18244	0.541	ng/ul#	91
29) Benzo(g,h,i)perylene	27.22	276	31436	1.069	ng/ul	98

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

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