

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121420\
 Data File : BM028115.D
 Acq On : 14 Dec 2020 16:18
 Operator : JU/CG
 Sample : L4958-04DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2S38DL

Quant Time: Dec 14 16:58:32 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	2226	0.40	ng/ul	0.00
4) Naphthalene-d8	11.05	136	8768	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.84	164	4992	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.56	188	10872	0.40	ng/ul	0.00
17) Chrysene-d12	21.67	240	9734	0.40	ng/ul	0.00
23) Perylene-d12	24.05	264	8854	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.43	96	779	0.31	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.62	152	2353	0.18	ng/ul	0.00
18) Fluoranthene-d10	19.56	212	4875	0.17	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.10	128	7889	0.319	ng/ul	99
7) 2-Methylnaphthalene	12.69	142	3704	0.227	ng/ul	98
12) Fluorene	15.88	166	6086	0.302	ng/ul	99
14) Pentachlorophenol	17.21	266	4071	1.475	ng/ul	99
16) Anthracene	17.69	178	11860	0.405	ng/ul	99
19) Fluoranthene	19.58	202	12147	0.317	ng/ul	97
20) Pyrene	19.95	202	13905	0.364	ng/ul	99
22) Chrysene	21.71	228	9871	0.286	ng/ul	99
24) Benzo(b)fluoranthene	23.32	252	8948	0.251	ng/ul	92
26) Benzo(a)pyrene	23.94	252	10195	0.346	ng/ul	97
28) Dibenzo(a,h)anthracene	26.49	278	19898	0.665	ng/ul	97

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

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